



wwPDB EM Validation Summary Report ⓘ

Dec 19, 2022 – 03:14 pm GMT

PDB ID : 7NRC
EMDB ID : EMD-12534
Title : Structure of the yeast Gcn1 bound to a leading stalled 80S ribosome with Rbg2, Gir2, A- and P-tRNA and eIF5A
Authors : Pochopien, A.A.; Beckert, B.; Wilson, D.N.
Deposited on : 2021-03-03
Resolution : 3.90 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

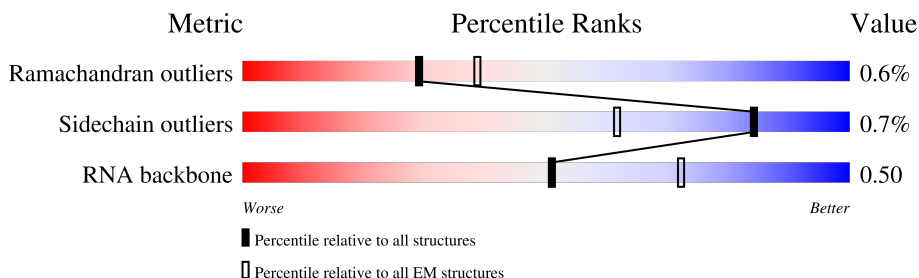
EMDB validation analysis : 0.0.1.dev43
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	S2	1800	
2	SI	6	
3	SP	206	
4	SQ	232	
5	SE	117	
6	SR	216	
7	SA	222	
8	SS	258	

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Mol	Chain	Length	Quality of chain
9	SB	206	9% 99%
10	ST	228	98%
11	SU	184	14% 97%
12	SV	198	94% 6%
13	SW	184	8% 97%
14	SC	92	10% 99%
15	SX	142	99%
16	SD	121	26% 97%
17	SY	150	5% 98%
18	SZ	127	98%
19	SF	141	99%
20	SG	125	10% 97%
21	SH	145	7% 97%
22	SI	143	9% 99%
23	SJ	100	7% 99%
24	Sa	87	99%
25	Sb	129	98%
26	Sc	144	99%
27	Sd	134	99%
28	SK	82	20% 98%
29	Se	97	100%
30	Sf	81	9% 98%
31	SM	53	100%
32	Sg	57	7% 100%
33	SN	73	26% 96%

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Mol	Chain	Length	Quality of chain
34	SO	312	17% 99%
35	SL	63	6% 98%
36	Sm	76	43% 38% 18%
37	Sn	75	60% 32% 8%
38	So	366	24% 95% 5%
39	Sp	48	15% 100%
40	LA	3394	67% 24% 6%
41	LB	121	79% 20%
42	LC	158	73% 25%
43	LD	251	100%
44	LE	386	99%
45	LF	361	99%
46	LG	294	6% 99%
47	LH	175	94% 5%
48	LI	222	100%
49	LJ	233	98%
50	LK	191	99%
51	LL	218	5% 99%
52	LM	169	22% 98%
53	LN	193	98%
54	LO	136	100%
55	LP	203	99%
56	LQ	197	100%
57	LR	183	100%
58	LS	185	97%

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Mol	Chain	Length	Quality of chain
59	LT	188	98% 7%
60	LU	171	100%
61	LV	159	99%
62	LW	100	99%
63	LX	136	100%
64	LY	65	100%
65	LZ	121	100%
66	La	125	98%
67	Lb	135	99%
68	Lc	148	96%
69	Ld	58	95% 5% 5%
70	Le	96	96%
71	Lf	109	98% 6%
72	Lg	127	99%
73	Lh	106	98%
74	Li	112	100%
75	Lj	119	97%
76	Lk	99	99%
77	Ll	81	96%
78	Lm	77	100% 5%
79	Ln	50	100%
80	Lo	52	98%
81	Lp	25	100%
82	Lq	103	99%
83	Lr	91	95%

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Mol	Chain	Length	Quality of chain
84	Ls	154	
85	Lt	210	
86	A	1209	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
84	5CT	Ls	51	X	-	-	-

2 Entry composition i

There are 86 unique types of molecules in this entry. The entry contains 213576 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called 18S rRNA (1771-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	S2	1771	37739	16872	6683	12413	1771	0	0

- Molecule 2 is a RNA chain called RNA (5'-R(P*AP*UP*GP*AP*AP*A)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
2	SI	6	131	59	27	39	6	0	0

- Molecule 3 is a protein called 40S ribosomal protein S0-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	SP	206	1603	1030	284	287	2	0	0

- Molecule 4 is a protein called 40S ribosomal protein S1-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	SQ	226	1798	1139	330	325	4	0	0

- Molecule 5 is a protein called 40S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	SE	117	916	583	171	155	7	0	0

- Molecule 6 is a protein called 40S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	SR	216	1626	1042	287	295	2	0	0

- Molecule 7 is a protein called 40S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	SA	222	1729	1098	312	313	6	0	0

- Molecule 8 is a protein called 40S ribosomal protein S4-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	SS	258	2056	1308	387	358	3	0	0

- Molecule 9 is a protein called 40S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	SB	206	1605	1005	299	298	3	0	0

- Molecule 10 is a protein called 40S ribosomal protein S6-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	ST	228	1815	1138	351	323	3	0	0

- Molecule 11 is a protein called 40S ribosomal protein S7-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
11	SU	184	1473	946	263	264	0	0

- Molecule 12 is a protein called 40S ribosomal protein S8-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	SV	187	1476	916	295	263	2	0	0

- Molecule 13 is a protein called 40S ribosomal protein S9-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
13	SW	184	1479	935	285	258	1	0	0

- Molecule 14 is a protein called 40S ribosomal protein S10-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	SC	92	Total	C	N	O	S	0	0
			752	487	122	141	2		

- Molecule 15 is a protein called 40S ribosomal protein S11-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	SX	142	Total	C	N	O	S	0	0
			1142	733	217	189	3		

- Molecule 16 is a protein called 40S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	SD	121	Total	C	N	O	S	0	0
			875	551	153	169	2		

- Molecule 17 is a protein called 40S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	SY	150	Total	C	N	O	S	0	0
			1192	759	224	207	2		

- Molecule 18 is a protein called 40S ribosomal protein S14-B.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	SZ	127	Total	C	N	O	S	0	0
			923	568	185	167	3		

- Molecule 19 is a protein called 40S ribosomal protein S16-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
19	SF	141	Total	C	N	O	0	0
			1105	708	203	194		

- Molecule 20 is a protein called 40S ribosomal protein S17-B.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	SG	121	Total	C	N	O	S	0	0
			948	596	179	171	2		

- Molecule 21 is a protein called 40S ribosomal protein S18-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	SH	145	Total	C	N	O	S	0	0
			1188	741	237	208	2		

- Molecule 22 is a protein called 40S ribosomal protein S19-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	SI	143	Total	C	N	O	S	0	0
			1112	694	208	208	2		

- Molecule 23 is a protein called 40S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	SJ	100	Total	C	N	O	S	0	0
			797	506	144	146	1		

- Molecule 24 is a protein called 40S ribosomal protein S21-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	Sa	87	Total	C	N	O	S	0	0
			673	415	125	131	2		

- Molecule 25 is a protein called 40S ribosomal protein S22-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	Sb	129	Total	C	N	O	S	0	0
			1021	650	188	180	3		

- Molecule 26 is a protein called 40S ribosomal protein S23-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	Sc	144	Total	C	N	O	S	0	0
			1121	708	220	191	2		

- Molecule 27 is a protein called 40S ribosomal protein S24-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
27	Sd	134	Total	C	N	O	0	0
			1032	651	195	186		

- Molecule 28 is a protein called 40S ribosomal protein S25-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
28	SK	82	Total	C	N	O	0	0
			651	416	123	112		

- Molecule 29 is a protein called 40S ribosomal protein S26-B.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	Se	97	Total	C	N	O	S	0	0
			765	473	160	127	5		

- Molecule 30 is a protein called 40S ribosomal protein S27-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	Sf	81	Total	C	N	O	S	0	0
			610	382	110	113	5		

- Molecule 31 is a protein called 40S ribosomal protein S29-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	SM	53	Total	C	N	O	S	0	0
			442	274	92	72	4		

- Molecule 32 is a protein called 40S ribosomal protein S30-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	Sg	57	Total	C	N	O	S	0	0
			451	284	93	73	1		

- Molecule 33 is a protein called 40S ribosomal protein S31.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	SN	73	Total	C	N	O	S	0	0
			556	352	105	95	4		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
SN	97	ALA	LYS	variant	UNP P05759

- Molecule 34 is a protein called Guanine nucleotide-binding protein subunit beta-like protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	SO	312	Total	C	N	O	S	0	0
			2383	1514	409	452	8		

- Molecule 35 is a protein called 40S ribosomal protein S28-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	SL	63	Total	C	N	O	S	0	0
			491	303	96	91	1		

- Molecule 36 is a RNA chain called tRNA (76-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
36	Sm	76	Total	C	N	O	P	0	0
			1611	721	281	534	75		

- Molecule 37 is a RNA chain called tRNA (75-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
37	Sn	75	Total	C	N	O	P	0	0
			1606	716	297	518	75		

- Molecule 38 is a protein called Ribosome-interacting GTPase 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	So	347	Total	C	N	O	S	0	0
			2729	1723	478	518	10		

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
So	62	ALA	LYS	variant	UNP P53295
So	92	ALA	ILE	variant	UNP P53295
So	94	GLU	HIS	variant	UNP P53295
So	96	GLU	ALA	variant	UNP P53295
So	134	LYS	ARG	variant	UNP P53295
So	271	ALA	THR	variant	UNP P53295
So	299	ALA	TYR	variant	UNP P53295
So	303	ALA	ARG	variant	UNP P53295
So	318	ASP	ASN	variant	UNP P53295
So	361	GLU	ASP	variant	UNP P53295

- Molecule 39 is a protein called GIR2.

Mol	Chain	Residues	Atoms				AltConf	Trace
39	Sp	48	Total	C	N	O	0	0
			240	144	48	48		

- Molecule 40 is a RNA chain called 25S rRNA (3184-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
40	LA	3184	Total	C	N	O	P	0	0
			68091	30415	12259	22233	3184		

- Molecule 41 is a RNA chain called 5S rRNA (121-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
41	LB	121	Total	C	N	O	P	0	0
			2579	1152	461	845	121		

- Molecule 42 is a RNA chain called 5.8S rRNA (158-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
42	LC	158	Total	C	N	O	P	0	0
			3353	1500	586	1109	158		

- Molecule 43 is a protein called 60S ribosomal protein L2-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	LD	251	Total	C	N	O	S	0	0
			1899	1182	385	331	1		

- Molecule 44 is a protein called 60S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	LE	386	Total	C	N	O	S	0	0
			3075	1950	584	533	8		

- Molecule 45 is a protein called 60S ribosomal protein L4-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	LF	361	Total	C	N	O	S	0	0
			2748	1729	522	494	3		

- Molecule 46 is a protein called 60S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
46	LG	294	2351	1484	410	455	2	0	0

- Molecule 47 is a protein called 60S ribosomal protein L6-B.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
47	LH	167	1307	843	234	230		0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
LH	120	LYS	ASN	variant	UNP P05739

- Molecule 48 is a protein called 60S ribosomal protein L7-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
48	LI	222	1784	1151	324	308	1	0	0

- Molecule 49 is a protein called 60S ribosomal protein L8-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
49	LJ	233	1804	1151	323	327	3	0	0

- Molecule 50 is a protein called 60S ribosomal protein L9-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
50	LK	191	1508	957	274	273	4	0	0

- Molecule 51 is a protein called 60S ribosomal protein L10.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
51	LL	218	1764	1117	334	306	7	0	0

- Molecule 52 is a protein called 60S ribosomal protein L11-B.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
52	LM	169	1346	843	252	247	4	0	0

- Molecule 53 is a protein called 60S ribosomal protein L13-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
53	LN	193	1543	962	315	266		0	0

- Molecule 54 is a protein called 60S ribosomal protein L14-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
54	LO	136	1053	675	199	177	2	0	0

- Molecule 55 is a protein called 60S ribosomal protein L15-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
55	LP	203	1720	1077	361	281	1	0	0

- Molecule 56 is a protein called 60S ribosomal protein L16-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
56	LQ	197	1555	1003	289	262	1	197	0

- Molecule 57 is a protein called 60S ribosomal protein L17-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
57	LR	183	1416	879	284	253		0	0

- Molecule 58 is a protein called 60S ribosomal protein L18-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
58	LS	185	1441	908	290	241	2	0	0

- Molecule 59 is a protein called 60S ribosomal protein L19-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
59	LT	188	1515	932	323	260	0	0

- Molecule 60 is a protein called 60S ribosomal protein L20-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
60	LU	171	1437	925	266	243	3	0	0

- Molecule 61 is a protein called 60S ribosomal protein L21-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
61	LV	159	1272	802	245	221	4	0	0

- Molecule 62 is a protein called 60S ribosomal protein L22-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
62	LW	100	796	516	131	149	0	0

- Molecule 63 is a protein called 60S ribosomal protein L23-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
63	LX	136	1003	628	189	179	7	0	0

- Molecule 64 is a protein called 60S ribosomal protein L24-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
64	LY	65	528	339	104	84	1	0	0

- Molecule 65 is a protein called 60S ribosomal protein L25.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
65	LZ	121	964	620	169	173	2	0	0

- Molecule 66 is a protein called 60S ribosomal protein L26-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
66	La	125	984	620	191	173	0	0

- Molecule 67 is a protein called 60S ribosomal protein L27-A.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
67	Lb	135	1080	701	199	180	0	0

- Molecule 68 is a protein called 60S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
68	Lc	148	1169	747	231	188	3	0	0

- Molecule 69 is a protein called 60S ribosomal protein L29.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
69	Ld	58	462	289	100	73	0	0

- Molecule 70 is a protein called 60S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
70	Le	96	737	476	123	137	1	0	0

- Molecule 71 is a protein called 60S ribosomal protein L31-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
71	Lf	109	876	556	167	152	1	0	0

- Molecule 72 is a protein called 60S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
72	Lg	127	1017	644	205	167	1	0	0

- Molecule 73 is a protein called 60S ribosomal protein L33-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
73	Lh	106	850	540	165	144	1	0	0

- Molecule 74 is a protein called 60S ribosomal protein L34-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
74	Li	112	880	545	179	152	4	0	0

- Molecule 75 is a protein called 60S ribosomal protein L35-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
75	Lj	119	969	615	186	167	1	0	0

- Molecule 76 is a protein called 60S ribosomal protein L36-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
76	Lk	99	766	478	154	132	2	0	0

- Molecule 77 is a protein called 60S ribosomal protein L37-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
77	Ll	81	645	393	141	106	5	0	0

- Molecule 78 is a protein called 60S ribosomal protein L38.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
78	Lm	77	612	391	115	106	0	0

- Molecule 79 is a protein called 60S ribosomal protein L39.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
79	Ln	50	436	272	97	65	2	0	0

- Molecule 80 is a protein called 60S ribosomal protein L40-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
80	Lo	52	Total	C	N	O	S	0	0
			410	254	86	65	5		

- Molecule 81 is a protein called 60S ribosomal protein L41-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
81	Lp	25	Total	C	N	O	S	0	0
			229	139	62	27	1		

- Molecule 82 is a protein called 60S ribosomal protein L42-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
82	Lq	103	Total	C	N	O	S	0	0
			824	517	167	135	5		

- Molecule 83 is a protein called 60S ribosomal protein L43-A.

Mol	Chain	Residues	Atoms					AltConf	Trace
83	Lr	91	Total	C	N	O	S	0	0
			694	429	138	121	6		

- Molecule 84 is a protein called eiF5A.

Mol	Chain	Residues	Atoms					AltConf	Trace
84	Ls	154	Total	C	N	O	S	0	0
			1143	709	195	230	9		

- Molecule 85 is a protein called L1 60S ribosomal protein.

Mol	Chain	Residues	Atoms				AltConf	Trace
85	Lt	210	Total	C	N	O	0	0
			1050	630	210	210		

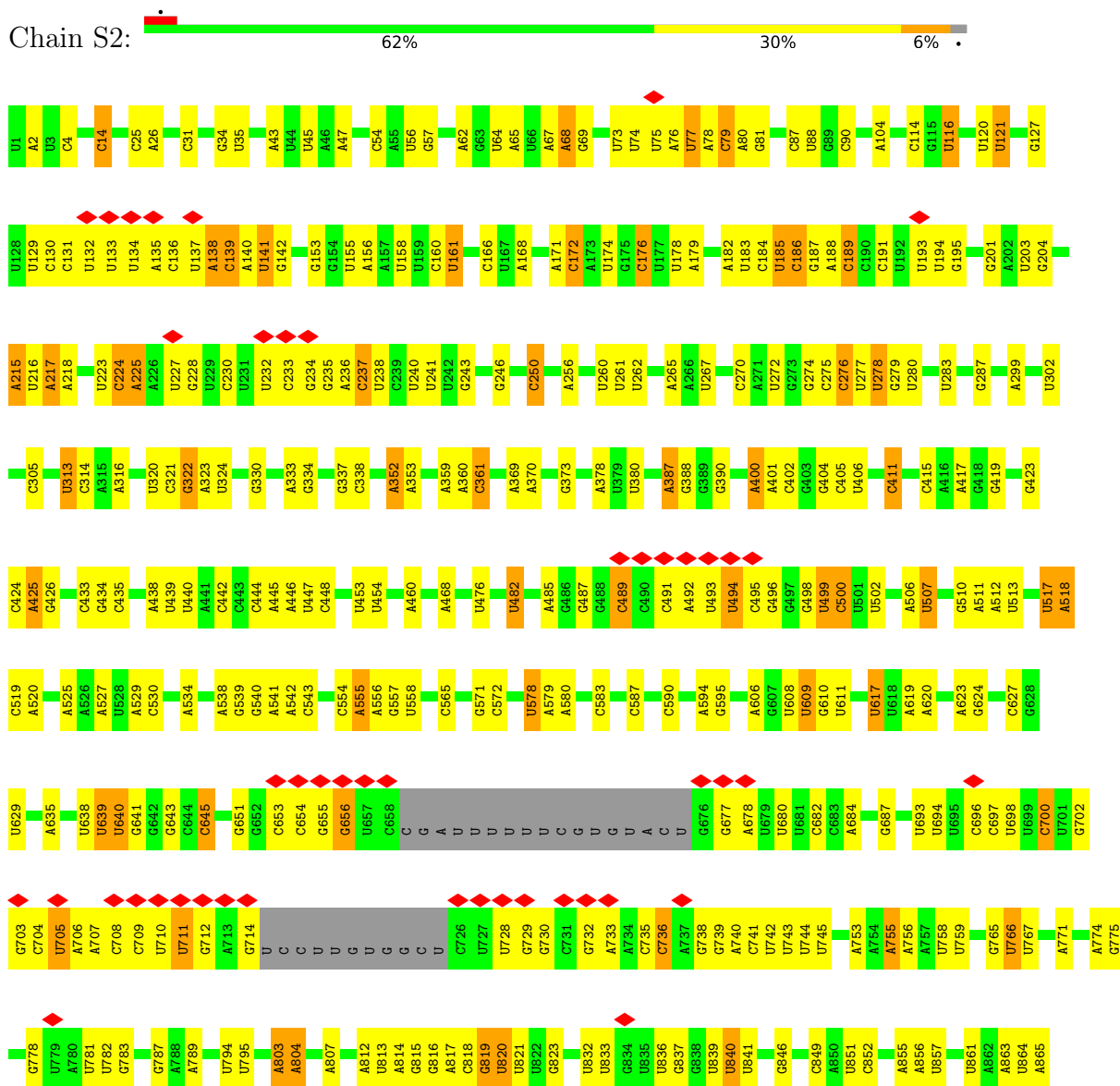
- Molecule 86 is a protein called GCN1.

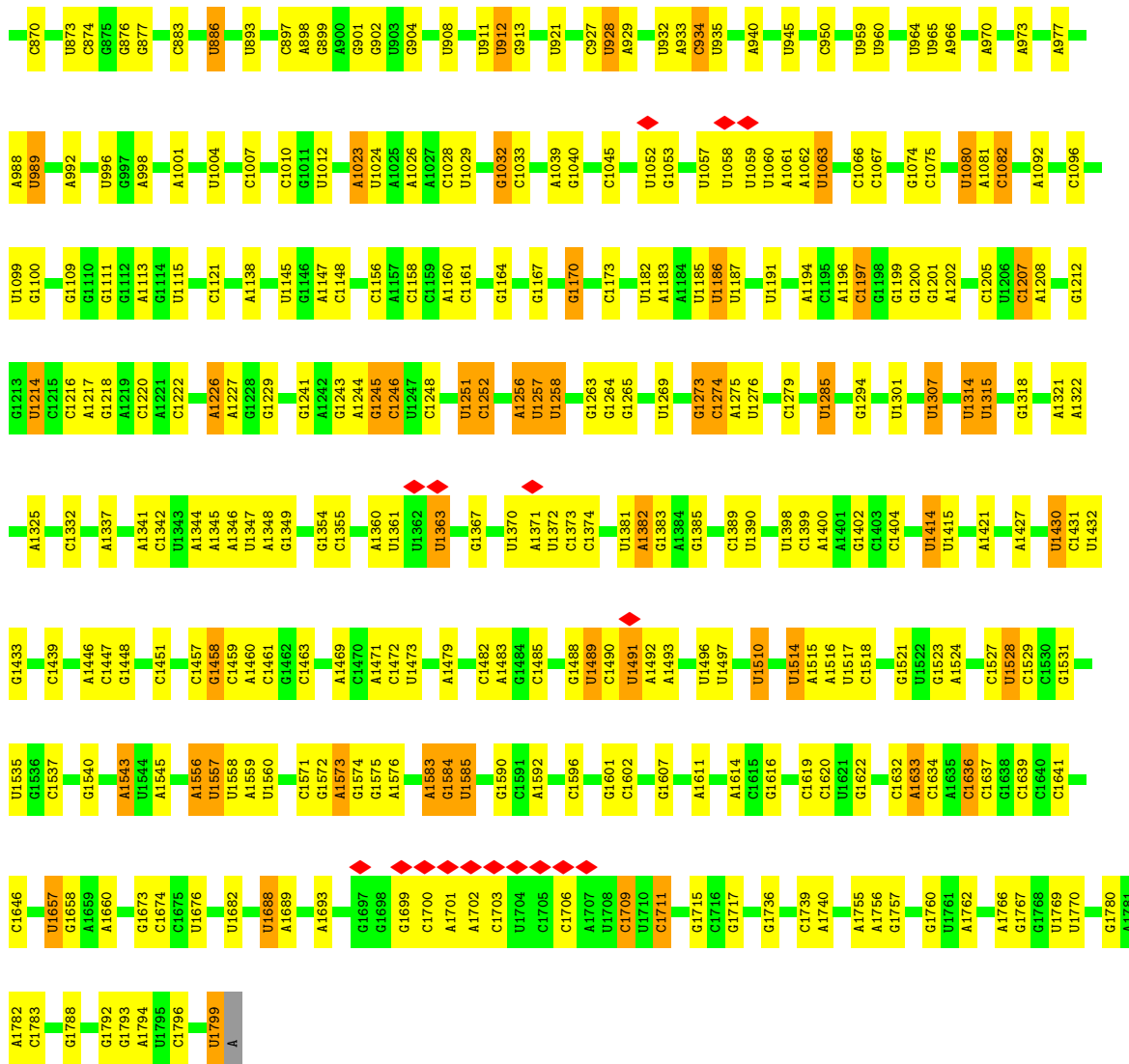
Mol	Chain	Residues	Atoms				AltConf	Trace
86	A	1209	Total	C	N	O	1	0
			6029	3609	1210	1210		

3 Residue-property plots i

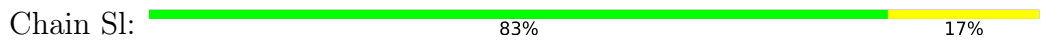
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: 18S rRNA (1771-MER)

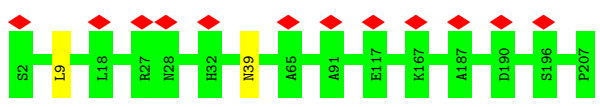




- Molecule 2: RNA (5'-R(P*AP*UP*GP*AP*AP*A)-3')

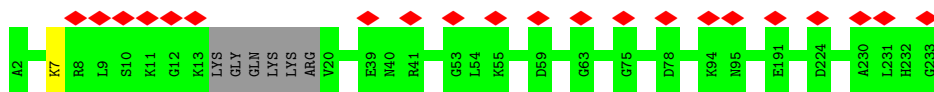


- Molecule 3: 40S ribosomal protein S0-A

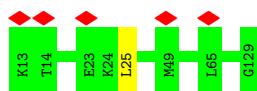


- Molecule 4: 40S ribosomal protein S1-A

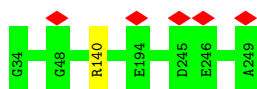




- Molecule 5: 40S ribosomal protein S15



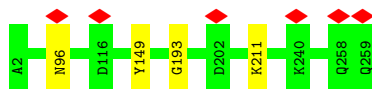
- Molecule 6: 40S ribosomal protein S2



- Molecule 7: 40S ribosomal protein S3



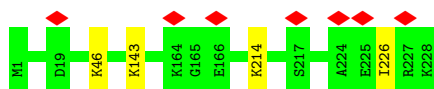
- Molecule 8: 40S ribosomal protein S4-A



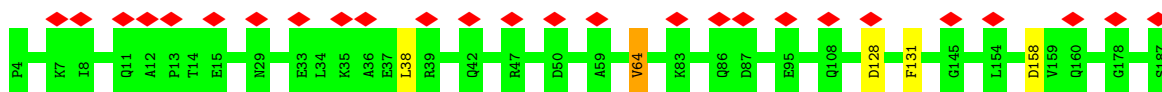
- Molecule 9: 40S ribosomal protein S5



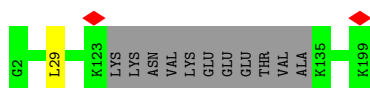
- Molecule 10: 40S ribosomal protein S6-A



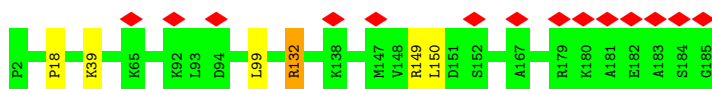
- Molecule 11: 40S ribosomal protein S7-A



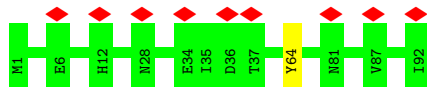
- Molecule 12: 40S ribosomal protein S8-A



- Molecule 13: 40S ribosomal protein S9-A



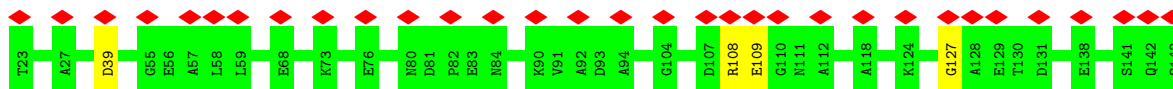
- Molecule 14: 40S ribosomal protein S10-A



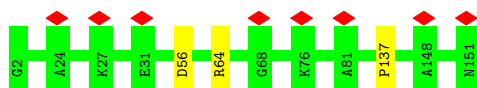
- Molecule 15: 40S ribosomal protein S11-A



- Molecule 16: 40S ribosomal protein S12

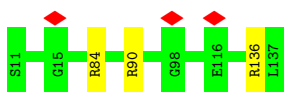


- Molecule 17: 40S ribosomal protein S13



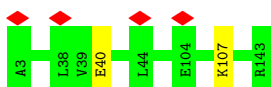
- Molecule 18: 40S ribosomal protein S14-B

Chain SZ:  98%



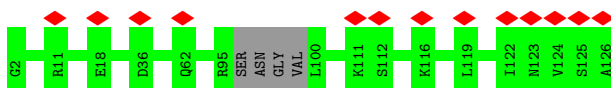
- Molecule 19: 40S ribosomal protein S16-A

Chain SF:  99%



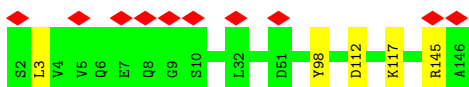
- Molecule 20: 40S ribosomal protein S17-B

Chain SG:  97%



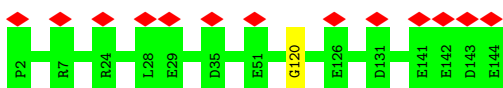
- Molecule 21: 40S ribosomal protein S18-A

Chain SH:  97%



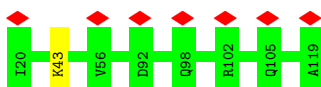
- Molecule 22: 40S ribosomal protein S19-A

Chain SI:  99%



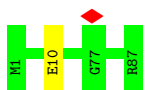
- Molecule 23: 40S ribosomal protein S20

Chain SJ:  99%



- Molecule 24: 40S ribosomal protein S21-A

Chain Sa:  99%



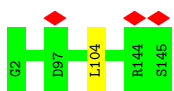
- Molecule 25: 40S ribosomal protein S22-A

Chain Sb: 98%



- Molecule 26: 40S ribosomal protein S23-A

Chain Sc: 99%



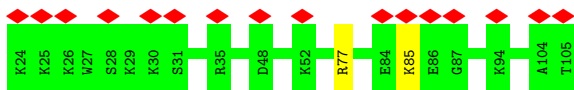
- Molecule 27: 40S ribosomal protein S24-A

Chain Sd: 99%



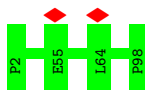
- Molecule 28: 40S ribosomal protein S25-A

Chain SK: 20% 98%



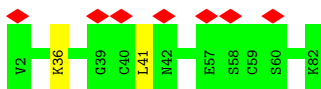
- Molecule 29: 40S ribosomal protein S26-B

Chain Se: 100%



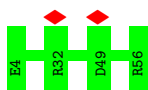
- Molecule 30: 40S ribosomal protein S27-A

Chain Sf: 9% 98%



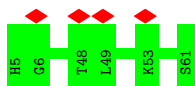
- Molecule 31: 40S ribosomal protein S29-A

Chain SM:  100%



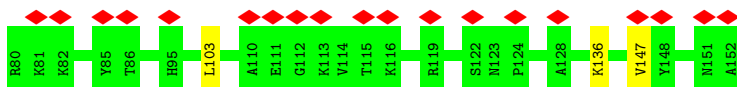
- Molecule 32: 40S ribosomal protein S30-A

Chain Sg:  7%



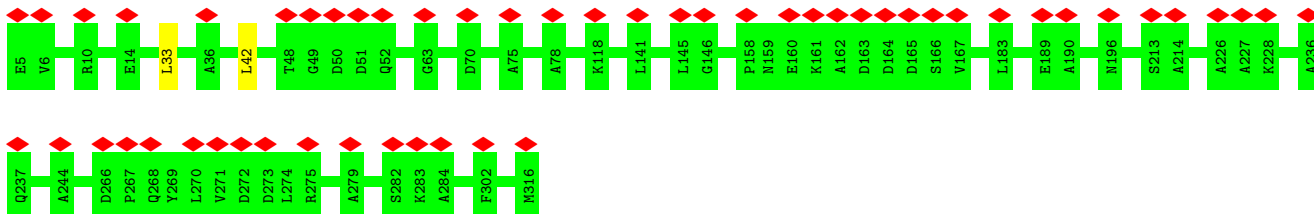
- Molecule 33: 40S ribosomal protein S31

Chain SN:  26%



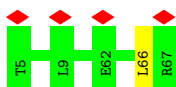
- Molecule 34: Guanine nucleotide-binding protein subunit beta-like protein

Chain SO:  17%



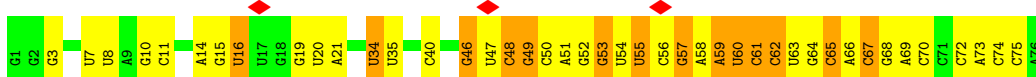
- Molecule 35: 40S ribosomal protein S28-A

Chain SL:  6%



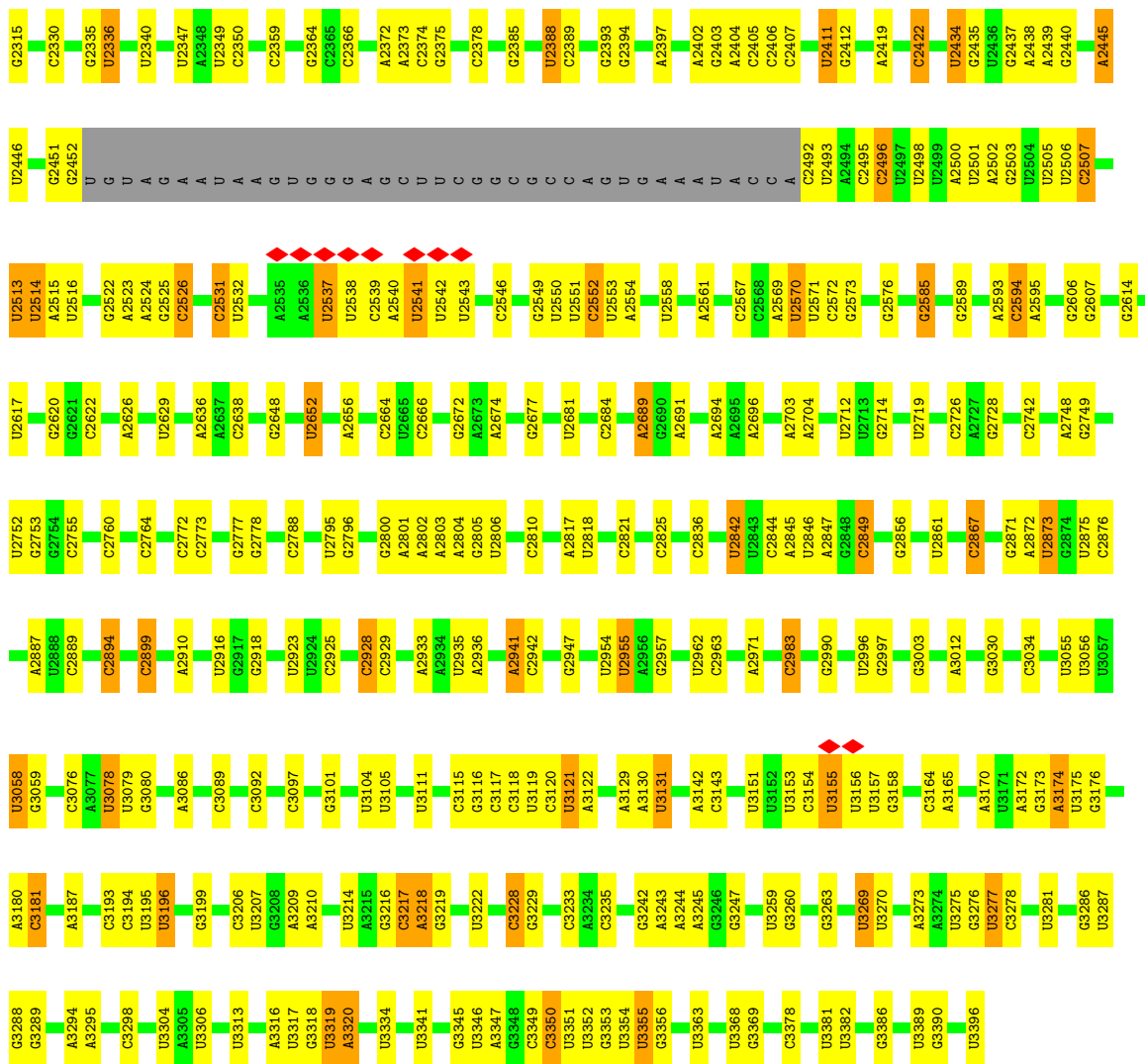
- Molecule 36: tRNA (76-MER)

Chain Sm:  43%

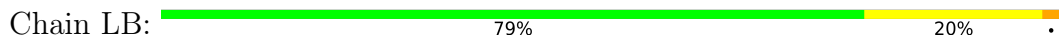


- Molecule 37: tRNA (75-MER)

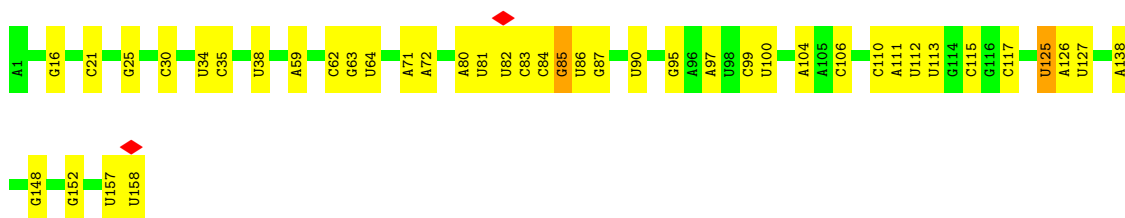
A621	A760	A917	C1032	A1190	C1272	A1419	U1540	A1642	A1797	C1951	A2187	G2187
G625	G761	A921	C1037	U1191	A1273	C1420	G1541	A1643	A1814	G1952	U1839	A2188
C637	G765	U922	C1040	C1192	A1274	G1421	G1542	U1645	U1815	G1953	A1840	A2189
C638	A766	C923	U1041	A1193	U1276	G1430	U1549	U1646	A1816	G1954	C1817	G2194
U642	G800	U932	C1045	C1196	C1277	G1431	U1555	G1658	U1818	U1955	U1819	G2201
A649	C803	A933	A1046	A1197	A1278	C1432	C1556	C1660	U1820	A	U1820	U2205
C650	A1047	G934	A1047	C1201	C1280	A1433	A1557	G1677	U1821	A	C1821	G2206
A650	A1048	G937	A1048	A1202	G1281	G1434	A1558	A1683	C1822	A	C1822	A2207
A660	A806	C944	C1049	A1206	G1282	C1437	G1560	C1685	C1832	A	C1832	A2208
C663	A808	C949	U1052	G1207	C1284	U1442	C1562	C1686	A	A	C1839	U2209
C667	A817	C949	A1064	U1208	G1285	G1443	U1563	U1686	A	A	U1840	A2223
U829	U829	U954	A1065	A1217	A1287	A1446	U1564	U1687	A	A	A1841	C2231
A830	A830	U954	A1065	U1218	A1287	G1447	G1565	U1688	A	A	A1842	A2232
G831	G831	C959	G1072	U1218	G1296	U1448	U1567	A1696	A	A	C1843	C2237
G832	G832	U960	C1076	G1222	C1296	A1449	U1568	A1704	A	A	G1845	A2244
A848	A848	C969	U1081	C1224	G1307	A1450	U1569	U1705	A	A	C1846	G2249
C849	C849	C971	U1081	C1227	A1308	U1455	U1570	C1706	A	A	C1849	G2250
U850	U850	C971	A1093	C1228	U1309	U1459	A1571	A1715	A	A	A1850	G2251
C851	C851	G974	U1094	G1229	G1313	U1470	U1572	U1716	A	A	C1854	A2252
U855	U855	C977	U1095	G1230	G1316	U1481	C1574	U1717	A	A	C1866	C2101
G860	G860	C977	U1096	U1234	A1317	A1482	G1575	U1722	A	A	A1867	U2102
C861	C861	C979	A1098	U1235	A1318	G1483	G1576	A1723	A	A	C1872	C2102
G712	G712	U979	C1098	G1236	A1318	U1484	C1579	U1724	A	A	U1879	G2111
A716	A716	A980	A1103	G1237	U1325	U1484	U1580	C1725	A	A	A1874	U2112
A720	A720	C982	G1104	C1238	A1330	G1488	C1581	U1746	A	A	G1874	A2113
G726	G726	U985	G1117	C1239	A1333	G1488	A1582	U1751	A	A	G1878	G2121
G737	G737	C991	C1118	U1241	C1345	U1496	A1583	G1751	A	A	A1879	C2122
G742	G742	A992	U1128	G1242	U1348	C1499	U1587	C1756	A	A	U1880	A2131
C743	C743	C993	G1145	G1243	G1349	C1499	A1588	U1759	A	A	A1881	C2132
C753	C753	G993	C1146	A1244	A1350	A1503	A1589	A1760	A	A	G1889	U2133
C758	C758	A1002	G1147	A1245	U1351	C1508	G1590	C1761	A	A	A1893	G2137
U764	U764	U1015	U1154	U1246	A1352	U1511	C1597	U1762	A	A	A1897	A2138
C765	C765	C1016	A1153	G1247	U1353	C1516	A1602	U1763	A	A	A1901	A2139
U766	U766	C1017	A1154	U1248	G1354	C1516	A1603	U1764	A	A	G1906	C2146
U767	U767	C1018	A1154	U1258	A1355	U1523	G1604	U1765	A	A	C1907	A2149
A771	A771	G1018	C1155	U1258	G1356	A1524	A1605	G1766	A	A	A1908	G2150
G774	G774	U1020	U1158	G1262	C1358	G1525	C1608	G1770	A	A	U1920	C2151
A775	A775	U1022	A1159	A1263	U1368	G1528	C1609	G1775	A	A	A1920	A2158
U776	U776	C1023	A1159	G1264	U1386	U1531	U1621	G1779	A	A	A1930	C2163
U777	U777	G1024	A1177	U1267	A1392	C1531	U1629	G1780	A	A	G1985	G2169
A914	A914	A1025	G1178	U1267	G1392	U1533	U1632	C1788	A	A	C1943	U2170
U776	U776	A1026	A1179	G1268	A1399	U1533	A1632	U1795	A	A	C1943	U2176
A915	A915	U1027	A1180	U1269	A1400	G1536	A1639	G1796	A	A	C1943	U2314
U1028	U1028	A1027	A1180	U1270	A1418	G1536	C1639	G1796	A	A	C1943	U2314
G1029	G1029	U1029	A1182	A1271	A1418	G1536	C1639	G1796	A	A	C1943	U2314



• Molecule 41: 5S rRNA (121-MER)



• Molecule 42: 5.8S rRNA (158-MER)



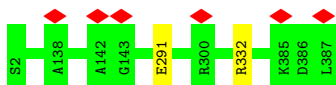
- Molecule 43: 60S ribosomal protein L2-A

Chain LD:  100%



- Molecule 44: 60S ribosomal protein L3

Chain LE:  99%



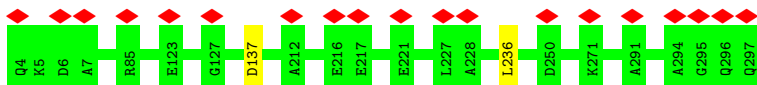
- Molecule 45: 60S ribosomal protein L4-A

Chain LF:  99%



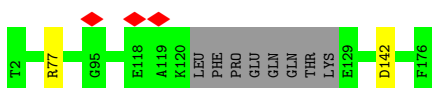
- Molecule 46: 60S ribosomal protein L5

Chain LG:  6%  99%



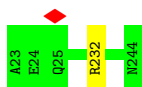
- Molecule 47: 60S ribosomal protein L6-B

Chain LH:  94%  5%



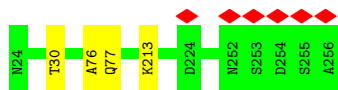
- Molecule 48: 60S ribosomal protein L7-A

Chain LI:  100%

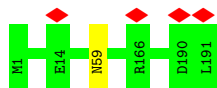


- Molecule 49: 60S ribosomal protein L8-A

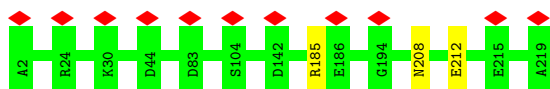
Chain LJ:  98%



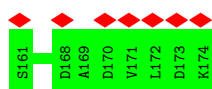
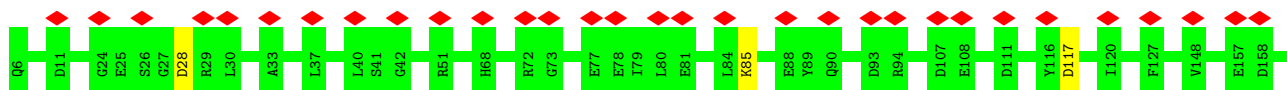
- Molecule 50: 60S ribosomal protein L9-A



- Molecule 51: 60S ribosomal protein L10



- Molecule 52: 60S ribosomal protein L11-B



- Molecule 53: 60S ribosomal protein L13-A



- Molecule 54: 60S ribosomal protein L14-A



There are no outlier residues recorded for this chain.

- Molecule 55: 60S ribosomal protein L15-A



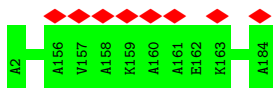
- Molecule 56: 60S ribosomal protein L16-A

Chain LQ:  100%

There are no outlier residues recorded for this chain.

- Molecule 57: 60S ribosomal protein L17-A

Chain LR:  100%



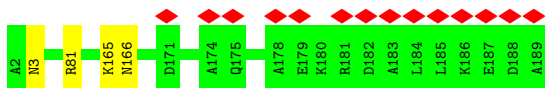
- Molecule 58: 60S ribosomal protein L18-A

Chain LS:  97%



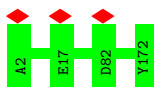
- Molecule 59: 60S ribosomal protein L19-A

Chain LT:  98%



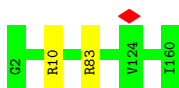
- Molecule 60: 60S ribosomal protein L20-A

Chain LU:  100%



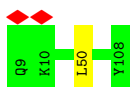
- Molecule 61: 60S ribosomal protein L21-A

Chain LV:  99%

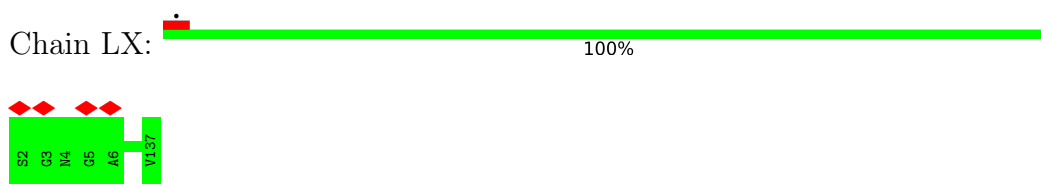


- Molecule 62: 60S ribosomal protein L22-A

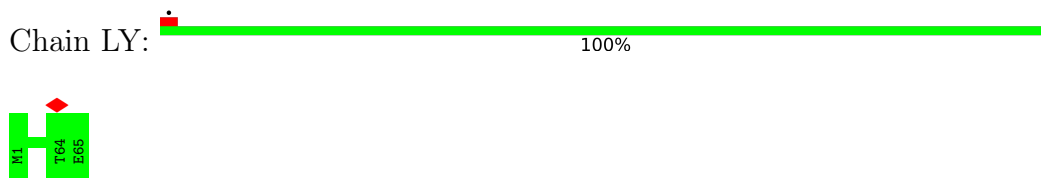
Chain LW:  99%



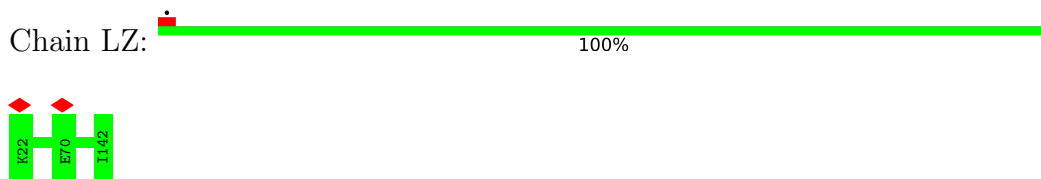
- Molecule 63: 60S ribosomal protein L23-A



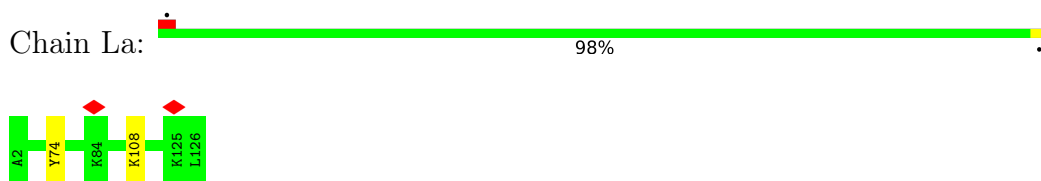
- Molecule 64: 60S ribosomal protein L24-A



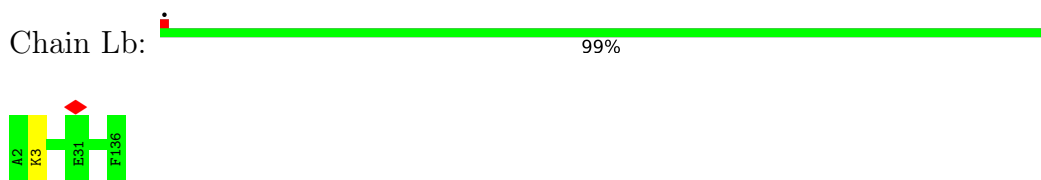
- Molecule 65: 60S ribosomal protein L25



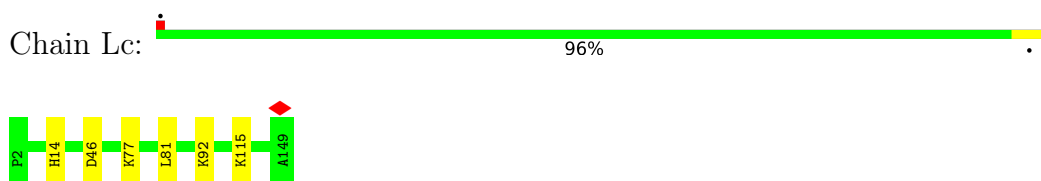
- Molecule 66: 60S ribosomal protein L26-A



- Molecule 67: 60S ribosomal protein L27-A

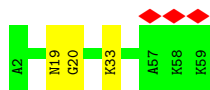


- Molecule 68: 60S ribosomal protein L28



- Molecule 69: 60S ribosomal protein L29

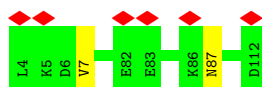




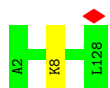
- Molecule 70: 60S ribosomal protein L30



- Molecule 71: 60S ribosomal protein L31-A



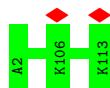
- Molecule 72: 60S ribosomal protein L32



- Molecule 73: 60S ribosomal protein L33-A



- Molecule 74: 60S ribosomal protein L34-A

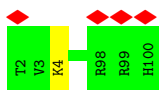


- Molecule 75: 60S ribosomal protein L35-A



- Molecule 76: 60S ribosomal protein L36-A

Chain Lk:  99%



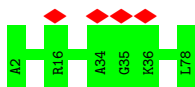
- Molecule 77: 60S ribosomal protein L37-A

Chain Ll:  96%



- Molecule 78: 60S ribosomal protein L38

Chain Lm:  100%



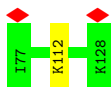
- Molecule 79: 60S ribosomal protein L39

Chain Ln:  100%

There are no outlier residues recorded for this chain.

- Molecule 80: 60S ribosomal protein L40-A

Chain Lo:  98%



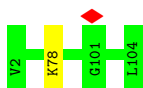
- Molecule 81: 60S ribosomal protein L41-A

Chain Lp:  100%

There are no outlier residues recorded for this chain.

- Molecule 82: 60S ribosomal protein L42-A

Chain Lq:  99%



- Molecule 83: 60S ribosomal protein L43-A

Chain Lr:  95%



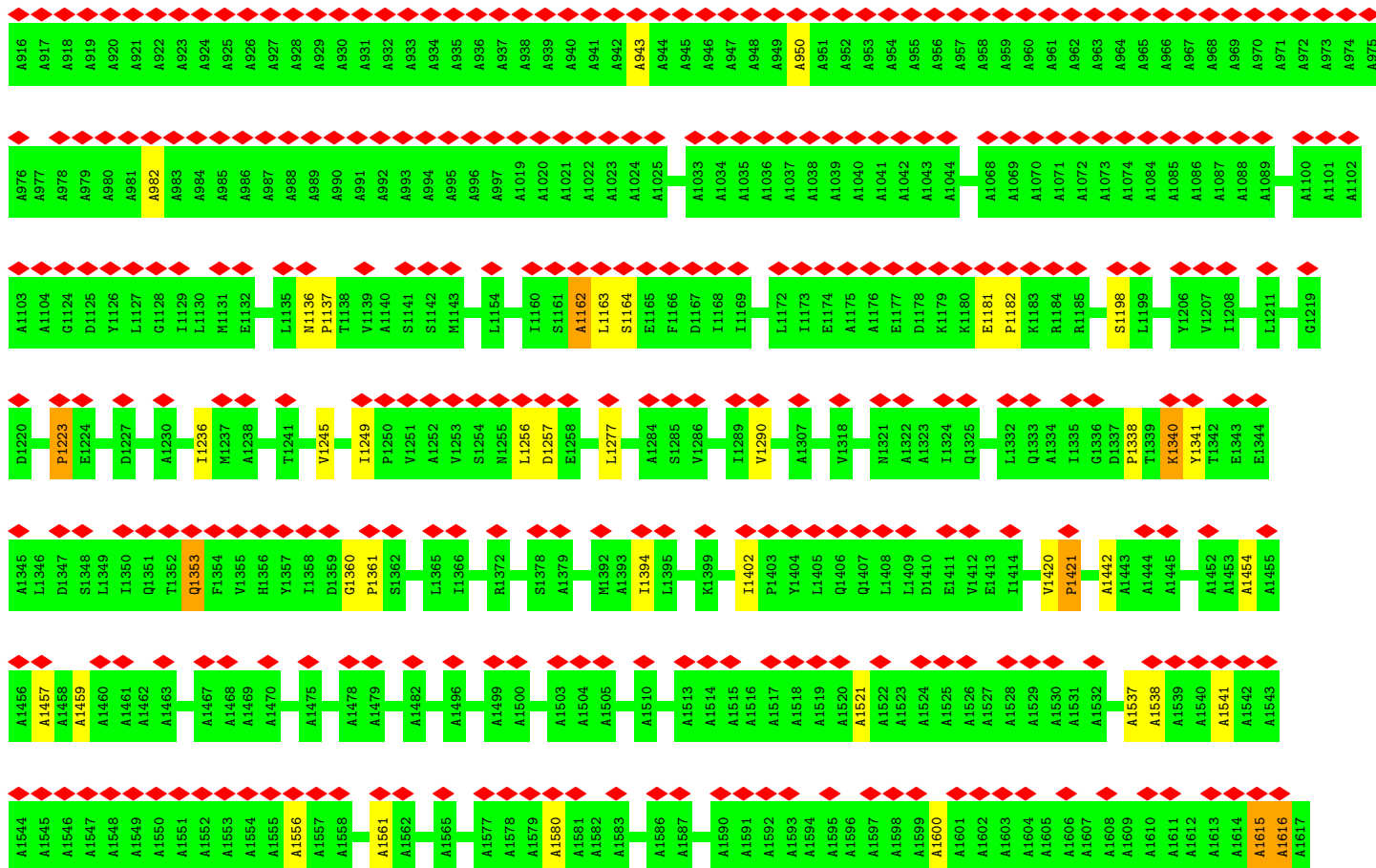
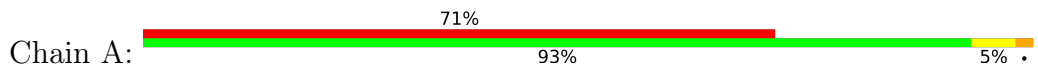
• Molecule 84: eiF5A

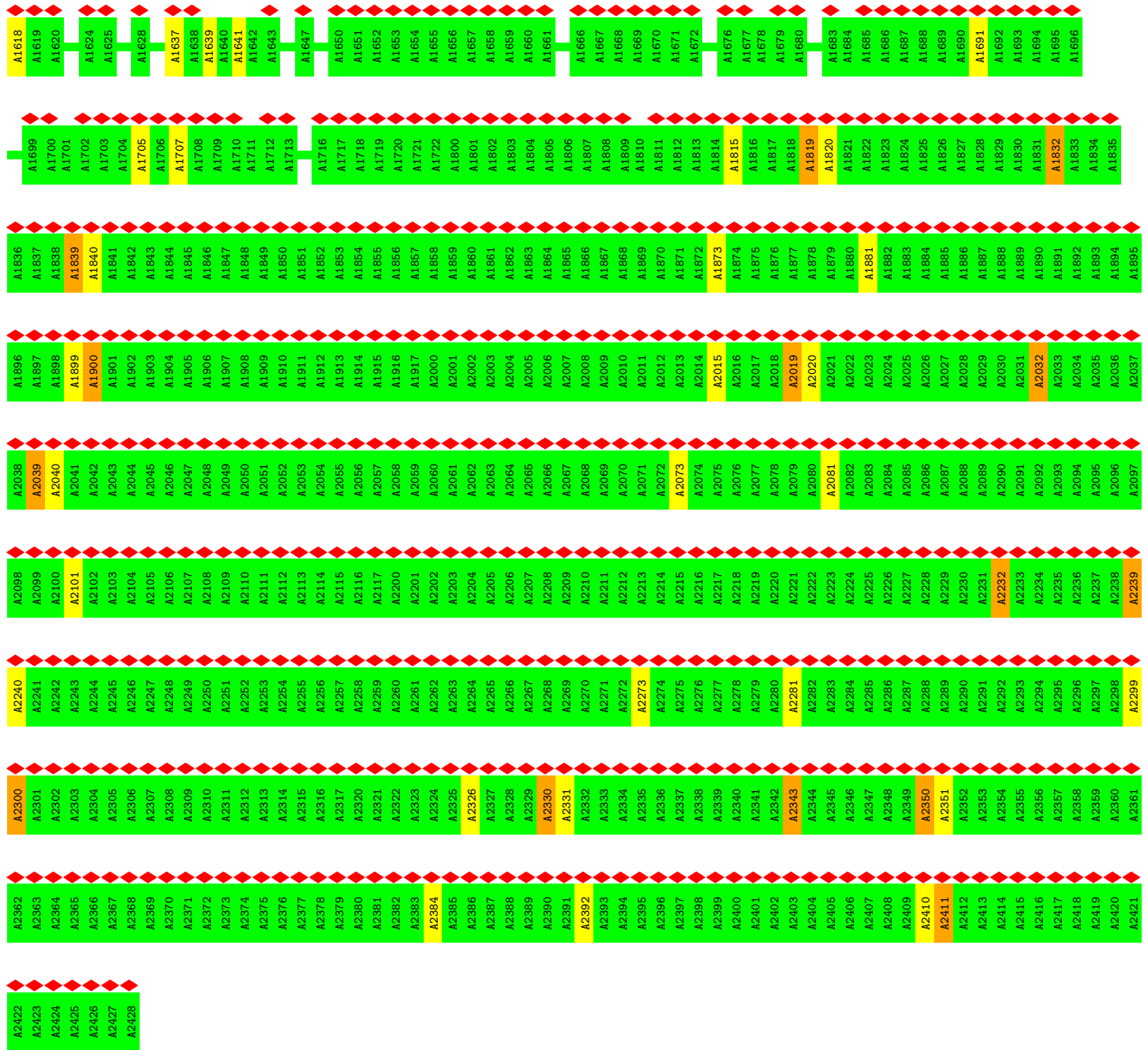


• Molecule 85: L1 60S ribosomal protein



• Molecule 86: GCN1





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	30016	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	2.5	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	0.512	Depositor
Minimum map value	-0.389	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.03	Depositor
Map size (Å)	758.8, 758.8, 758.8	wwPDB
Map dimensions	700, 700, 700	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.084, 1.084, 1.084	Depositor

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: 5CT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	S2	0.46	0/42211	1.27	410/65773 (0.6%)
2	SI	0.52	0/147	1.18	0/227
3	SP	0.37	0/1644	0.74	1/2249 (0.0%)
4	SQ	0.37	0/1823	0.84	0/2447
5	SE	0.37	0/936	0.80	1/1259 (0.1%)
6	SR	0.40	0/1656	0.81	2/2251 (0.1%)
7	SA	0.36	0/1754	0.76	0/2361
8	SS	0.39	0/2097	0.84	2/2823 (0.1%)
9	SB	0.36	0/1625	0.73	0/2197
10	ST	0.38	0/1839	0.78	0/2460
11	SU	0.40	0/1498	0.86	5/2019 (0.2%)
12	SV	0.35	0/1501	0.75	1/2006 (0.0%)
13	SW	0.33	0/1504	0.76	3/2016 (0.1%)
14	SC	0.37	0/769	0.78	0/1039
15	SX	0.34	0/1168	0.73	0/1575
16	SD	0.39	0/883	0.87	1/1199 (0.1%)
17	SY	0.38	0/1215	0.78	2/1638 (0.1%)
18	SZ	0.37	0/934	0.82	2/1257 (0.2%)
19	SF	0.37	0/1125	0.79	0/1510
20	SG	0.34	0/957	0.74	0/1283
21	SH	0.36	0/1207	0.80	3/1623 (0.2%)
22	SI	0.39	0/1130	0.76	0/1517
23	SJ	0.35	0/807	0.77	0/1091
24	Sa	0.33	0/682	0.75	1/921 (0.1%)
25	Sb	0.39	0/1038	0.83	2/1395 (0.1%)
26	Sc	0.39	0/1139	0.78	1/1518 (0.1%)
27	Sd	0.38	0/1046	0.79	2/1401 (0.1%)
28	SK	0.39	0/661	0.89	2/888 (0.2%)
29	Se	0.35	0/778	0.74	0/1042
30	Sf	0.31	0/620	0.80	1/838 (0.1%)
31	SM	0.39	0/452	0.79	0/600
32	Sg	0.32	0/459	0.80	0/611

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	SN	0.37	0/567	0.85	2/764 (0.3%)
34	SO	0.34	0/2436	0.83	2/3318 (0.1%)
35	SL	0.39	0/493	0.95	1/663 (0.2%)
36	Sm	0.49	0/1799	1.53	43/2801 (1.5%)
37	Sn	0.46	0/1796	1.37	24/2799 (0.9%)
38	So	0.29	0/2769	0.57	0/3736
40	LA	0.47	1/76214 (0.0%)	1.20	582/118821 (0.5%)
41	LB	0.39	0/2883	1.15	20/4491 (0.4%)
42	LC	0.45	0/3746	1.19	25/5832 (0.4%)
43	LD	0.36	0/1933	0.76	0/2598
44	LE	0.38	0/3146	0.74	1/4228 (0.0%)
45	LF	0.36	0/2800	0.73	0/3790
46	LG	0.37	0/2400	0.74	2/3239 (0.1%)
47	LH	0.37	0/1329	0.76	2/1794 (0.1%)
48	LI	0.38	0/1821	0.68	0/2451
49	LJ	0.37	0/1836	0.72	0/2481
50	LK	0.36	0/1529	0.74	0/2060
51	LL	0.35	0/1801	0.73	1/2416 (0.0%)
52	LM	0.36	0/1367	0.80	2/1834 (0.1%)
53	LN	0.36	0/1568	0.71	1/2106 (0.0%)
54	LO	0.35	0/1068	0.69	0/1438
55	LP	0.44	0/1757	0.75	1/2354 (0.0%)
56	LQ	0.39	0/1585	0.69	0/2128
57	LR	0.37	0/1439	0.69	0/1938
58	LS	0.35	0/1465	0.72	1/1965 (0.1%)
59	LT	0.34	0/1532	0.65	0/2043
60	LU	0.36	0/1473	0.69	0/1980
61	LV	0.35	0/1296	0.71	0/1739
62	LW	0.37	0/812	0.73	1/1099 (0.1%)
63	LX	0.34	0/1018	0.74	0/1369
64	LY	0.31	0/540	0.56	0/717
65	LZ	0.36	0/979	0.74	0/1321
66	La	0.31	0/995	0.68	0/1329
67	Lb	0.33	0/1106	0.68	0/1485
68	Lc	0.36	0/1200	0.74	2/1607 (0.1%)
69	Ld	0.33	0/473	0.69	0/629
70	Le	0.36	0/745	0.76	1/1001 (0.1%)
71	Lf	0.36	0/890	0.75	0/1196
72	Lg	0.33	0/1038	0.69	0/1390
73	Lh	0.37	0/868	0.81	1/1168 (0.1%)
74	Li	0.40	0/890	0.72	0/1189
75	Lj	0.32	0/978	0.70	1/1301 (0.1%)
76	Lk	0.34	0/772	0.69	0/1026

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
77	Ll	0.45	0/660	0.81	0/875
78	Lm	0.36	0/618	0.77	0/826
79	Ln	0.34	0/443	0.74	0/588
80	Lo	0.33	0/416	0.72	0/553
81	Lp	0.37	0/230	0.74	0/296
82	Lq	0.37	0/836	0.76	0/1104
83	Lr	0.38	0/701	0.72	1/934 (0.1%)
84	Ls	0.39	0/1142	0.99	10/1537 (0.7%)
86	A	0.92	21/6015 (0.3%)	0.89	19/8386 (0.2%)
All	All	0.45	22/227518 (0.0%)	1.07	1188/333767 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
8	SS	0	1
9	SB	0	2
11	SU	0	1
16	SD	0	1
18	SZ	0	1
19	SF	0	1
22	SI	0	1
25	Sb	0	1
33	SN	0	1
45	LF	0	3
48	LI	0	1
49	LJ	0	3
58	LS	0	1
68	Lc	0	2
69	Ld	0	2
70	Le	0	1
71	Lf	0	1
75	Lj	0	1
84	Ls	1	2
85	Lt	0	4
86	A	0	12
All	All	1	43

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
86	A	1361	PRO	N-CA	13.63	1.70	1.47
86	A	1421	PRO	N-CA	13.12	1.69	1.47
86	A	1223	PRO	N-CA	13.01	1.69	1.47
86	A	2351	ALA	N-CA	12.07	1.70	1.46
86	A	1840	ALA	N-CA	12.04	1.70	1.46

The worst 5 of 1188 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
40	LA	3217	C	N1-C2-O2	13.82	127.19	118.90
37	Sn	49	C	C6-N1-C2	-12.90	115.14	120.30
40	LA	2492	C	N1-C2-O2	12.72	126.53	118.90
86	A	1340	LYS	O-C-N	-12.26	103.08	122.70
86	A	2019	ALA	O-C-N	-12.22	103.15	122.70

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
84	Ls	51	5CT	C2

5 of 43 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
9	SB	125	THR	Peptide
9	SB	42	LEU	Peptide
16	SD	108	ARG	Peptide
8	SS	193	GLY	Peptide
11	SU	64	VAL	Peptide

5.2 Too-close contacts [i](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	SP	204/206 (99%)	183 (90%)	21 (10%)	0	100	100
4	SQ	222/232 (96%)	201 (90%)	21 (10%)	0	100	100
5	SE	115/117 (98%)	100 (87%)	15 (13%)	0	100	100
6	SR	214/216 (99%)	190 (89%)	24 (11%)	0	100	100
7	SA	220/222 (99%)	209 (95%)	11 (5%)	0	100	100
8	SS	256/258 (99%)	229 (90%)	27 (10%)	0	100	100
9	SB	204/206 (99%)	189 (93%)	15 (7%)	0	100	100
10	ST	226/228 (99%)	211 (93%)	14 (6%)	1 (0%)	34	71
11	SU	182/184 (99%)	163 (90%)	19 (10%)	0	100	100
12	SV	183/198 (92%)	167 (91%)	16 (9%)	0	100	100
13	SW	182/184 (99%)	165 (91%)	16 (9%)	1 (0%)	29	67
14	SC	90/92 (98%)	79 (88%)	11 (12%)	0	100	100
15	SX	140/142 (99%)	124 (89%)	16 (11%)	0	100	100
16	SD	119/121 (98%)	91 (76%)	26 (22%)	2 (2%)	9	43
17	SY	148/150 (99%)	133 (90%)	15 (10%)	0	100	100
18	SZ	125/127 (98%)	110 (88%)	15 (12%)	0	100	100
19	SF	139/141 (99%)	125 (90%)	14 (10%)	0	100	100
20	SG	117/125 (94%)	112 (96%)	5 (4%)	0	100	100
21	SH	143/145 (99%)	131 (92%)	12 (8%)	0	100	100
22	SI	141/143 (99%)	128 (91%)	13 (9%)	0	100	100
23	SJ	98/100 (98%)	88 (90%)	10 (10%)	0	100	100
24	Sa	85/87 (98%)	73 (86%)	12 (14%)	0	100	100
25	Sb	127/129 (98%)	113 (89%)	14 (11%)	0	100	100
26	Sc	142/144 (99%)	123 (87%)	19 (13%)	0	100	100
27	Sd	132/134 (98%)	124 (94%)	8 (6%)	0	100	100
28	SK	80/82 (98%)	69 (86%)	11 (14%)	0	100	100
29	Se	95/97 (98%)	84 (88%)	11 (12%)	0	100	100
30	Sf	79/81 (98%)	67 (85%)	12 (15%)	0	100	100
31	SM	51/53 (96%)	48 (94%)	3 (6%)	0	100	100
32	Sg	55/57 (96%)	43 (78%)	12 (22%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
33	SN	71/73 (97%)	48 (68%)	23 (32%)	0	100	100
34	SO	310/312 (99%)	268 (86%)	42 (14%)	0	100	100
35	SL	61/63 (97%)	57 (93%)	4 (7%)	0	100	100
38	So	341/366 (93%)	311 (91%)	30 (9%)	0	100	100
43	LD	249/251 (99%)	223 (90%)	26 (10%)	0	100	100
44	LE	384/386 (100%)	355 (92%)	29 (8%)	0	100	100
45	LF	359/361 (99%)	323 (90%)	35 (10%)	1 (0%)	41	75
46	LG	292/294 (99%)	266 (91%)	26 (9%)	0	100	100
47	LH	163/175 (93%)	147 (90%)	16 (10%)	0	100	100
48	LI	220/222 (99%)	203 (92%)	17 (8%)	0	100	100
49	LJ	231/233 (99%)	208 (90%)	23 (10%)	0	100	100
50	LK	189/191 (99%)	174 (92%)	15 (8%)	0	100	100
51	LL	216/218 (99%)	189 (88%)	27 (12%)	0	100	100
52	LM	167/169 (99%)	153 (92%)	14 (8%)	0	100	100
53	LN	191/193 (99%)	171 (90%)	19 (10%)	1 (0%)	29	67
54	LO	134/136 (98%)	121 (90%)	13 (10%)	0	100	100
55	LP	201/203 (99%)	181 (90%)	19 (10%)	1 (0%)	29	67
56	LQ	195/197 (99%)	186 (95%)	9 (5%)	0	100	100
57	LR	181/183 (99%)	168 (93%)	13 (7%)	0	100	100
58	LS	183/185 (99%)	170 (93%)	13 (7%)	0	100	100
59	LT	186/188 (99%)	178 (96%)	8 (4%)	0	100	100
60	LU	169/171 (99%)	159 (94%)	10 (6%)	0	100	100
61	LV	157/159 (99%)	143 (91%)	14 (9%)	0	100	100
62	LW	98/100 (98%)	93 (95%)	5 (5%)	0	100	100
63	LX	134/136 (98%)	128 (96%)	6 (4%)	0	100	100
64	LY	63/65 (97%)	59 (94%)	4 (6%)	0	100	100
65	LZ	119/121 (98%)	110 (92%)	9 (8%)	0	100	100
66	La	123/125 (98%)	117 (95%)	6 (5%)	0	100	100
67	Lb	133/135 (98%)	117 (88%)	16 (12%)	0	100	100
68	Lc	146/148 (99%)	130 (89%)	16 (11%)	0	100	100
69	Ld	56/58 (97%)	47 (84%)	9 (16%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
70	Le	94/96 (98%)	92 (98%)	2 (2%)	0	100	100
71	Lf	107/109 (98%)	93 (87%)	14 (13%)	0	100	100
72	Lg	125/127 (98%)	115 (92%)	10 (8%)	0	100	100
73	Lh	104/106 (98%)	98 (94%)	6 (6%)	0	100	100
74	Li	110/112 (98%)	105 (96%)	5 (4%)	0	100	100
75	Lj	117/119 (98%)	109 (93%)	8 (7%)	0	100	100
76	Lk	97/99 (98%)	91 (94%)	6 (6%)	0	100	100
77	Ll	79/81 (98%)	71 (90%)	8 (10%)	0	100	100
78	Lm	75/77 (97%)	72 (96%)	3 (4%)	0	100	100
79	Ln	48/50 (96%)	46 (96%)	2 (4%)	0	100	100
80	Lo	50/52 (96%)	44 (88%)	6 (12%)	0	100	100
81	Lp	23/25 (92%)	23 (100%)	0	0	100	100
82	Lq	101/103 (98%)	93 (92%)	8 (8%)	0	100	100
83	Lr	89/91 (98%)	87 (98%)	2 (2%)	0	100	100
84	Ls	151/154 (98%)	117 (78%)	31 (20%)	3 (2%)	7	40
86	A	1182/1209 (98%)	929 (79%)	190 (16%)	63 (5%)	2	22
All	All	12588/12828 (98%)	11260 (90%)	1255 (10%)	73 (1%)	29	63

5 of 73 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
84	Ls	133	ASP
86	A	1162	ALA
86	A	1163	LEU
86	A	1164	SER
86	A	1182	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	SP	170/173 (98%)	169 (99%)	1 (1%)	86	91
4	SQ	200/205 (98%)	199 (100%)	1 (0%)	88	93
5	SE	95/98 (97%)	95 (100%)	0	100	100
6	SR	175/175 (100%)	175 (100%)	0	100	100
7	SA	182/182 (100%)	180 (99%)	2 (1%)	73	84
8	SS	220/220 (100%)	218 (99%)	2 (1%)	78	87
9	SB	172/173 (99%)	172 (100%)	0	100	100
10	ST	189/195 (97%)	186 (98%)	3 (2%)	62	79
11	SU	163/165 (99%)	163 (100%)	0	100	100
12	SV	148/159 (93%)	148 (100%)	0	100	100
13	SW	156/157 (99%)	153 (98%)	3 (2%)	57	75
14	SC	77/85 (91%)	76 (99%)	1 (1%)	69	82
15	SX	126/127 (99%)	124 (98%)	2 (2%)	62	79
16	SD	88/98 (90%)	88 (100%)	0	100	100
17	SY	127/127 (100%)	126 (99%)	1 (1%)	81	89
18	SZ	90/96 (94%)	89 (99%)	1 (1%)	73	84
19	SF	117/117 (100%)	116 (99%)	1 (1%)	78	87
20	SG	101/113 (89%)	101 (100%)	0	100	100
21	SH	127/128 (99%)	125 (98%)	2 (2%)	62	79
22	SI	115/115 (100%)	115 (100%)	0	100	100
23	SJ	93/93 (100%)	92 (99%)	1 (1%)	73	84
24	Sa	71/74 (96%)	71 (100%)	0	100	100
25	Sb	110/110 (100%)	110 (100%)	0	100	100
26	Sc	119/119 (100%)	119 (100%)	0	100	100
27	Sd	102/112 (91%)	101 (99%)	1 (1%)	76	86
28	SK	67/73 (92%)	66 (98%)	1 (2%)	65	80
29	Se	82/83 (99%)	82 (100%)	0	100	100
30	Sf	70/70 (100%)	69 (99%)	1 (1%)	67	81
31	SM	47/47 (100%)	47 (100%)	0	100	100
32	Sg	48/49 (98%)	48 (100%)	0	100	100
33	SN	56/63 (89%)	56 (100%)	0	100	100
34	SO	250/257 (97%)	250 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
35	SL	55/56 (98%)	55 (100%)	0	100	100
38	So	298/308 (97%)	298 (100%)	0	100	100
43	LD	190/193 (98%)	189 (100%)	1 (0%)	88	93
44	LE	318/322 (99%)	317 (100%)	1 (0%)	92	95
45	LF	288/288 (100%)	287 (100%)	1 (0%)	92	95
46	LG	241/243 (99%)	241 (100%)	0	100	100
47	LH	139/154 (90%)	139 (100%)	0	100	100
48	LI	186/186 (100%)	186 (100%)	0	100	100
49	LJ	187/191 (98%)	186 (100%)	1 (0%)	88	93
50	LK	168/171 (98%)	167 (99%)	1 (1%)	86	91
51	LL	185/185 (100%)	183 (99%)	2 (1%)	73	84
52	LM	145/147 (99%)	144 (99%)	1 (1%)	84	90
53	LN	154/154 (100%)	152 (99%)	2 (1%)	69	82
54	LO	107/107 (100%)	107 (100%)	0	100	100
55	LP	175/175 (100%)	175 (100%)	0	100	100
56	LQ	160/160 (100%)	160 (100%)	0	100	100
57	LR	138/145 (95%)	138 (100%)	0	100	100
58	LS	150/150 (100%)	146 (97%)	4 (3%)	44	67
59	LT	152/153 (99%)	148 (97%)	4 (3%)	46	68
60	LU	155/155 (100%)	155 (100%)	0	100	100
61	LV	135/136 (99%)	133 (98%)	2 (2%)	65	80
62	LW	87/87 (100%)	87 (100%)	0	100	100
63	LX	104/104 (100%)	104 (100%)	0	100	100
64	LY	54/57 (95%)	54 (100%)	0	100	100
65	LZ	104/105 (99%)	104 (100%)	0	100	100
66	La	108/108 (100%)	106 (98%)	2 (2%)	57	75
67	Lb	112/115 (97%)	111 (99%)	1 (1%)	78	87
68	Lc	117/118 (99%)	115 (98%)	2 (2%)	60	78
69	Ld	46/46 (100%)	45 (98%)	1 (2%)	52	71
70	Le	81/81 (100%)	79 (98%)	2 (2%)	47	69
71	Lf	92/96 (96%)	91 (99%)	1 (1%)	73	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
72	Lg	108/109 (99%)	107 (99%)	1 (1%)	78	87
73	Lh	90/90 (100%)	89 (99%)	1 (1%)	73	84
74	Li	95/95 (100%)	95 (100%)	0	100	100
75	Lj	104/104 (100%)	102 (98%)	2 (2%)	57	75
76	Lk	80/81 (99%)	79 (99%)	1 (1%)	69	82
77	Ll	67/67 (100%)	64 (96%)	3 (4%)	27	56
78	Lm	68/68 (100%)	68 (100%)	0	100	100
79	Ln	45/45 (100%)	45 (100%)	0	100	100
80	Lo	45/47 (96%)	44 (98%)	1 (2%)	52	71
81	Lp	22/23 (96%)	22 (100%)	0	100	100
82	Lq	87/88 (99%)	86 (99%)	1 (1%)	73	84
83	Lr	71/71 (100%)	66 (93%)	5 (7%)	15	44
84	Ls	118/129 (92%)	116 (98%)	2 (2%)	60	78
All	All	9584/9771 (98%)	9514 (99%)	70 (1%)	84	90

5 of 70 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
76	Lk	4	LYS
77	Ll	19	CYS
83	Lr	57	CYS
43	LD	242	ARG
30	Sf	36	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 66 such sidechains are listed below:

Mol	Chain	Res	Type
71	Lf	87	ASN
74	Li	52	GLN
84	Ls	88	ASN
27	Sd	34	ASN
25	Sb	80	ASN

5.3.3 RNA

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	S2	1768/1800 (98%)	533 (30%)	52 (2%)
2	Sl	5/6 (83%)	1 (20%)	0
36	Sm	75/76 (98%)	35 (46%)	0
37	Sn	74/75 (98%)	23 (31%)	0
40	LA	3180/3394 (93%)	716 (22%)	42 (1%)
41	LB	120/121 (99%)	15 (12%)	1 (0%)
42	LC	157/158 (99%)	32 (20%)	3 (1%)
All	All	5379/5630 (95%)	1355 (25%)	98 (1%)

5 of 1355 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	S2	2	A
1	S2	4	C
1	S2	14	C
1	S2	25	C
1	S2	26	A

5 of 98 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
40	LA	849	C
40	LA	1815	U
40	LA	896	A
40	LA	1307	G
40	LA	2445	A

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
84	5CT	Ls	51	84,40	13,14,15	2.31	1 (7%)	9,15,17	2.26	4 (44%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
84	5CT	Ls	51	84,40	1/1/2/4	5/13/14/16	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
84	Ls	51	5CT	O1-C2	-8.19	1.18	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
84	Ls	51	5CT	O1-C2-C3	3.62	119.57	109.21
84	Ls	51	5CT	O1-C2-C1	3.37	120.84	109.32
84	Ls	51	5CT	C3-C2-C1	3.26	119.58	112.16
84	Ls	51	5CT	C1-NZ-CE	-2.69	107.41	113.42

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
84	Ls	51	5CT	C2

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
84	Ls	51	5CT	C2-C1-NZ-CE
84	Ls	51	5CT	NZ-C1-C2-C3
84	Ls	51	5CT	O1-C2-C3-C4
84	Ls	51	5CT	C2-C3-C4-N1
84	Ls	51	5CT	CE-CD-CG-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
86	A	13

The worst 5 of 13 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	1104:ALA	C	1124:GLY	N	40.36
1	A	997:ALA	C	1019:ALA	N	35.44
1	A	1917:ALA	C	2000:ALA	N	14.67
1	A	1044:ALA	C	1055:ALA	N	12.76
1	A	1074:ALA	C	1084:ALA	N	11.98

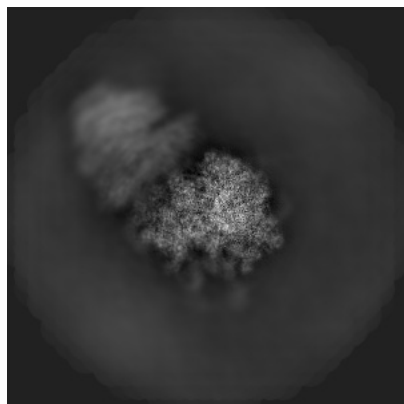
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-12534. These allow visual inspection of the internal detail of the map and identification of artifacts.

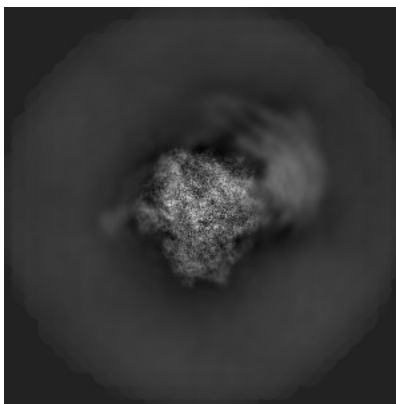
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

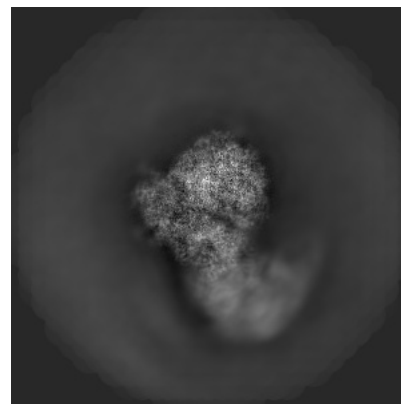
6.1.1 Primary map



X

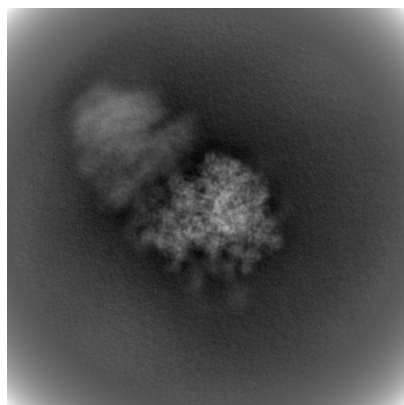


Y

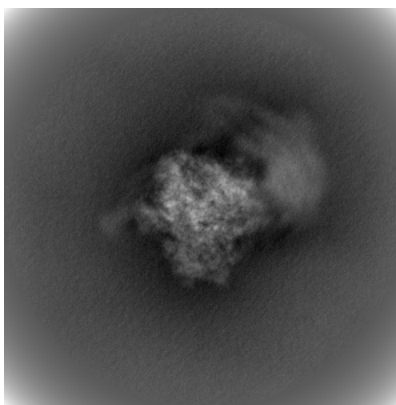


Z

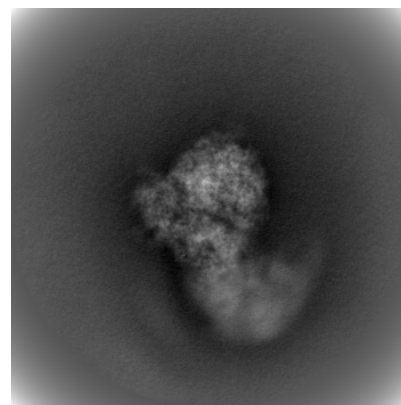
6.1.2 Raw map



X



Y

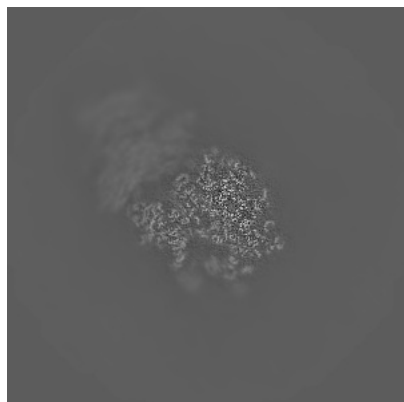


Z

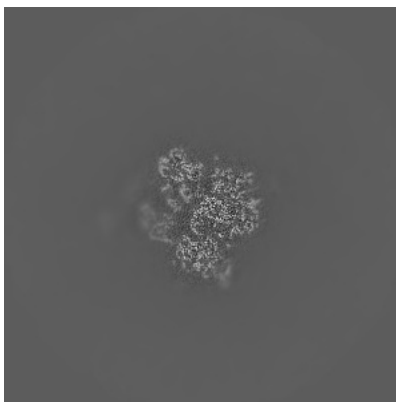
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

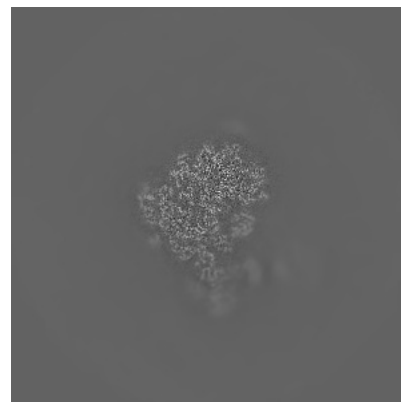
6.2.1 Primary map



X Index: 350

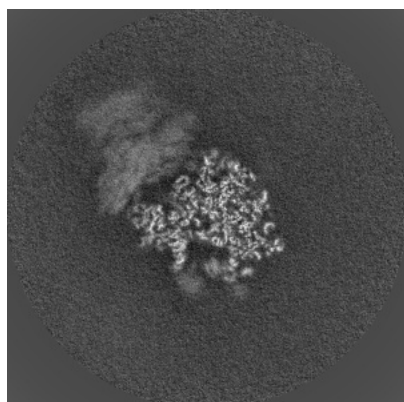


Y Index: 350

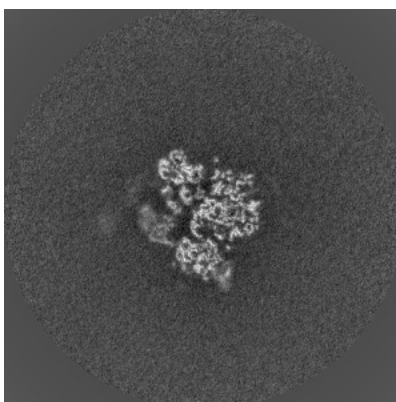


Z Index: 350

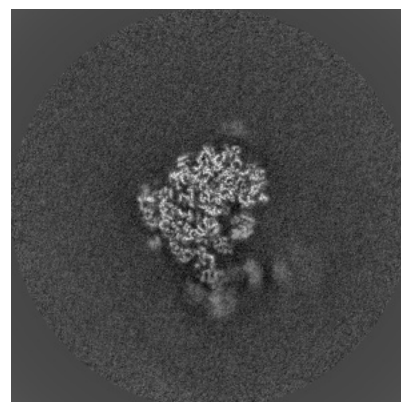
6.2.2 Raw map



X Index: 350



Y Index: 350

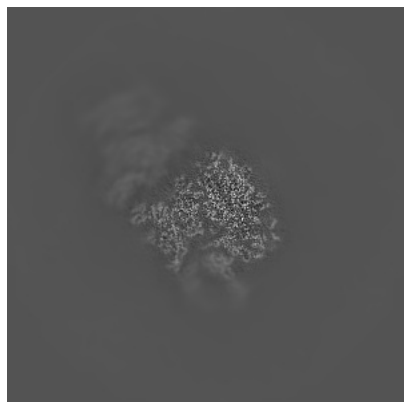


Z Index: 350

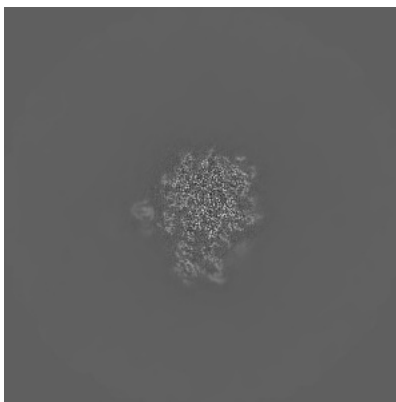
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 338

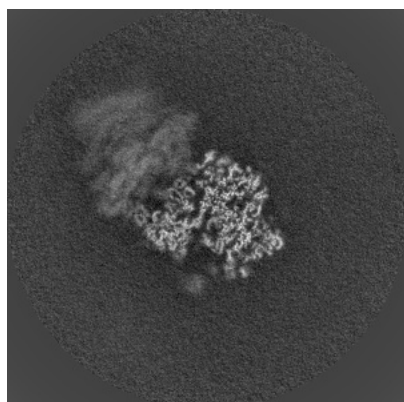


Y Index: 381

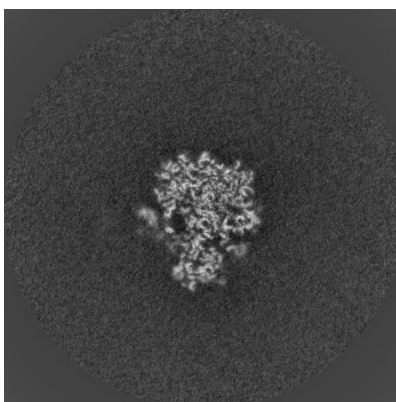


Z Index: 356

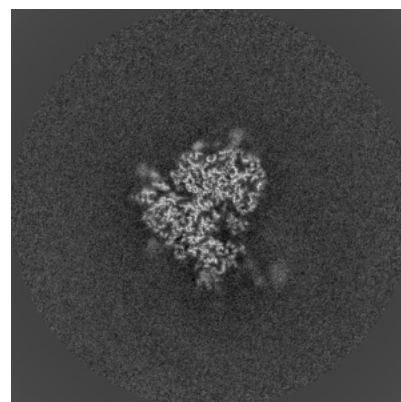
6.3.2 Raw map



X Index: 360



Y Index: 369



Z Index: 334

The images above show the largest variance slices of the map in three orthogonal directions.

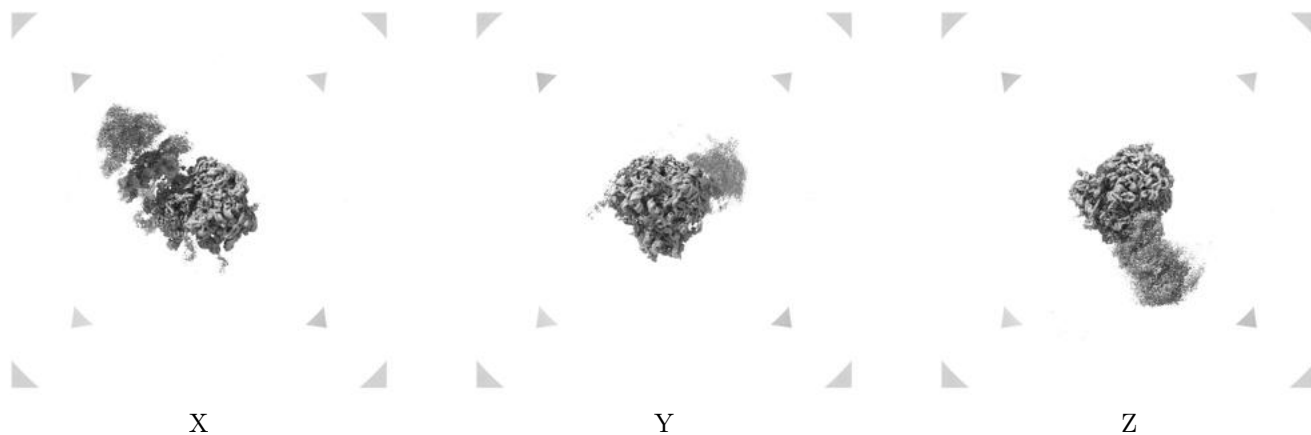
6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.03. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

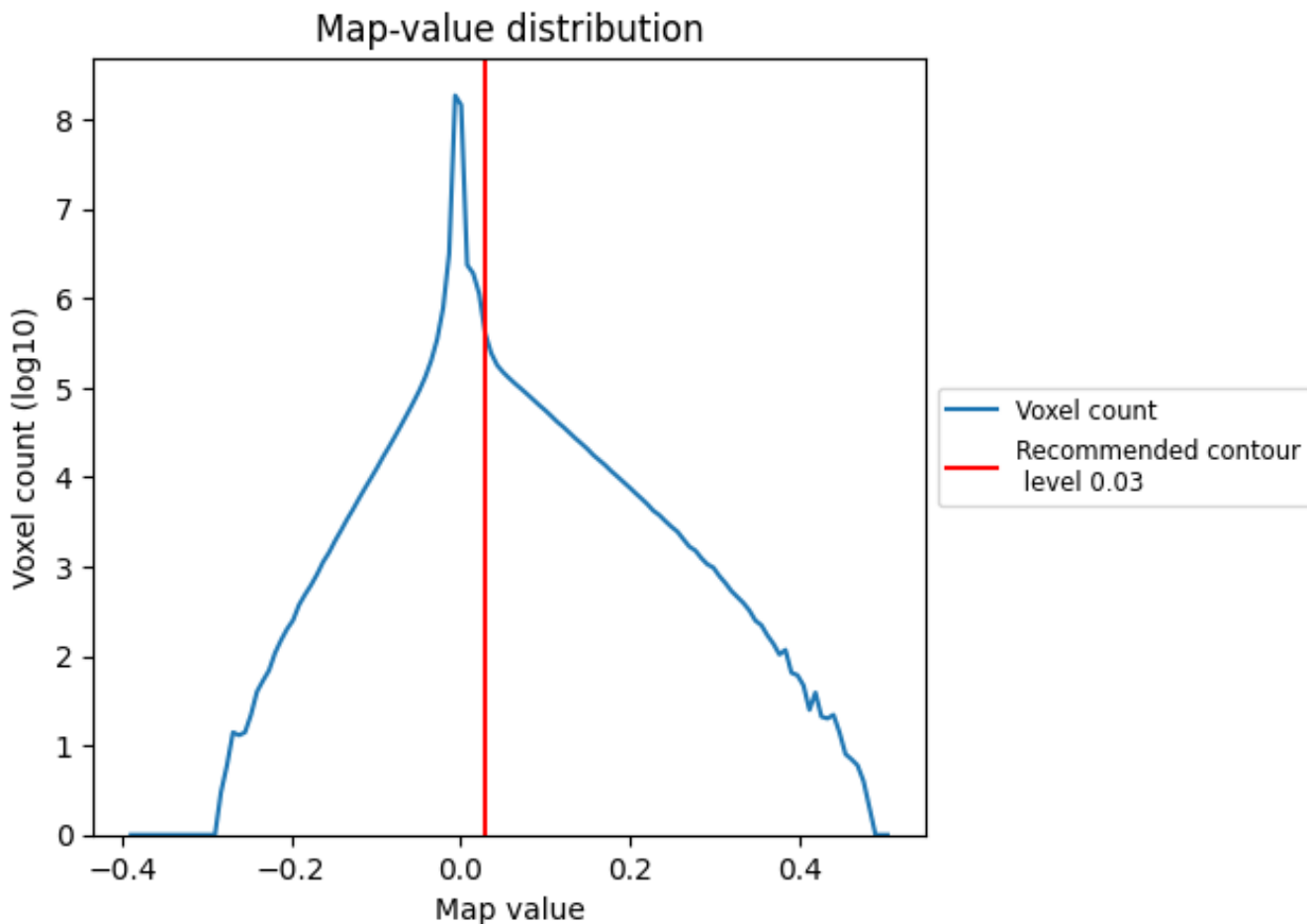
6.5 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

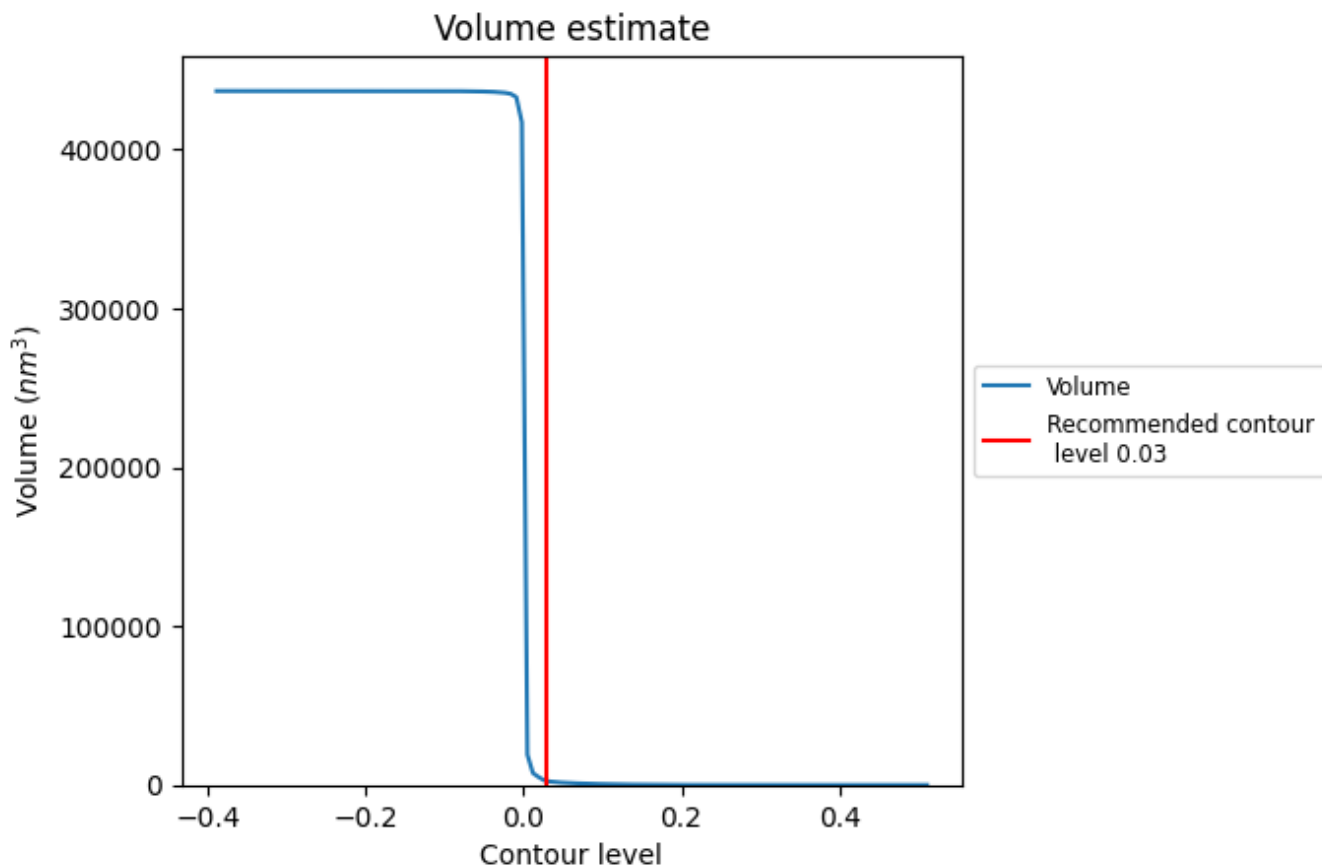
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

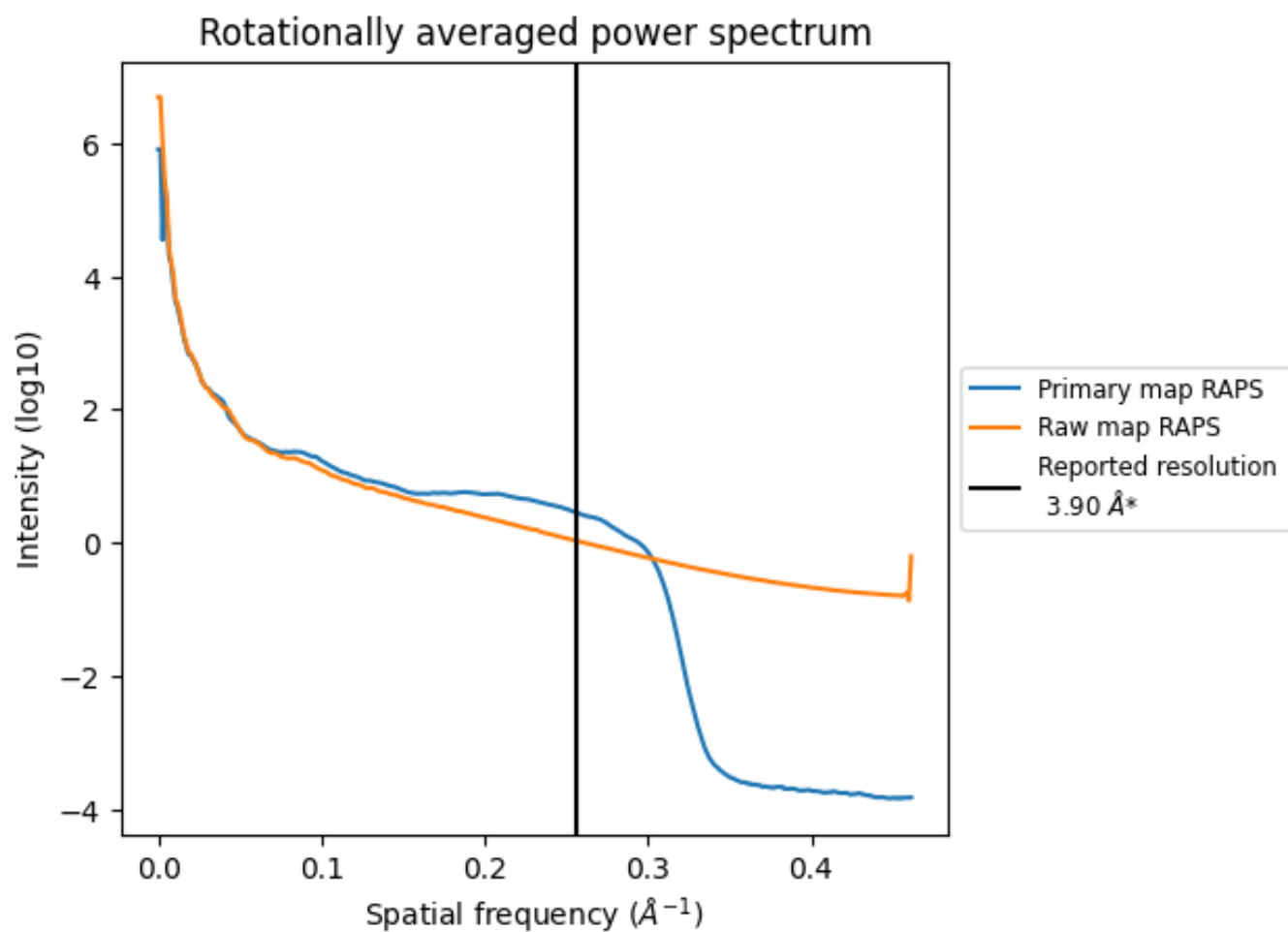
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2499 nm^3 ; this corresponds to an approximate mass of 2257 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum i

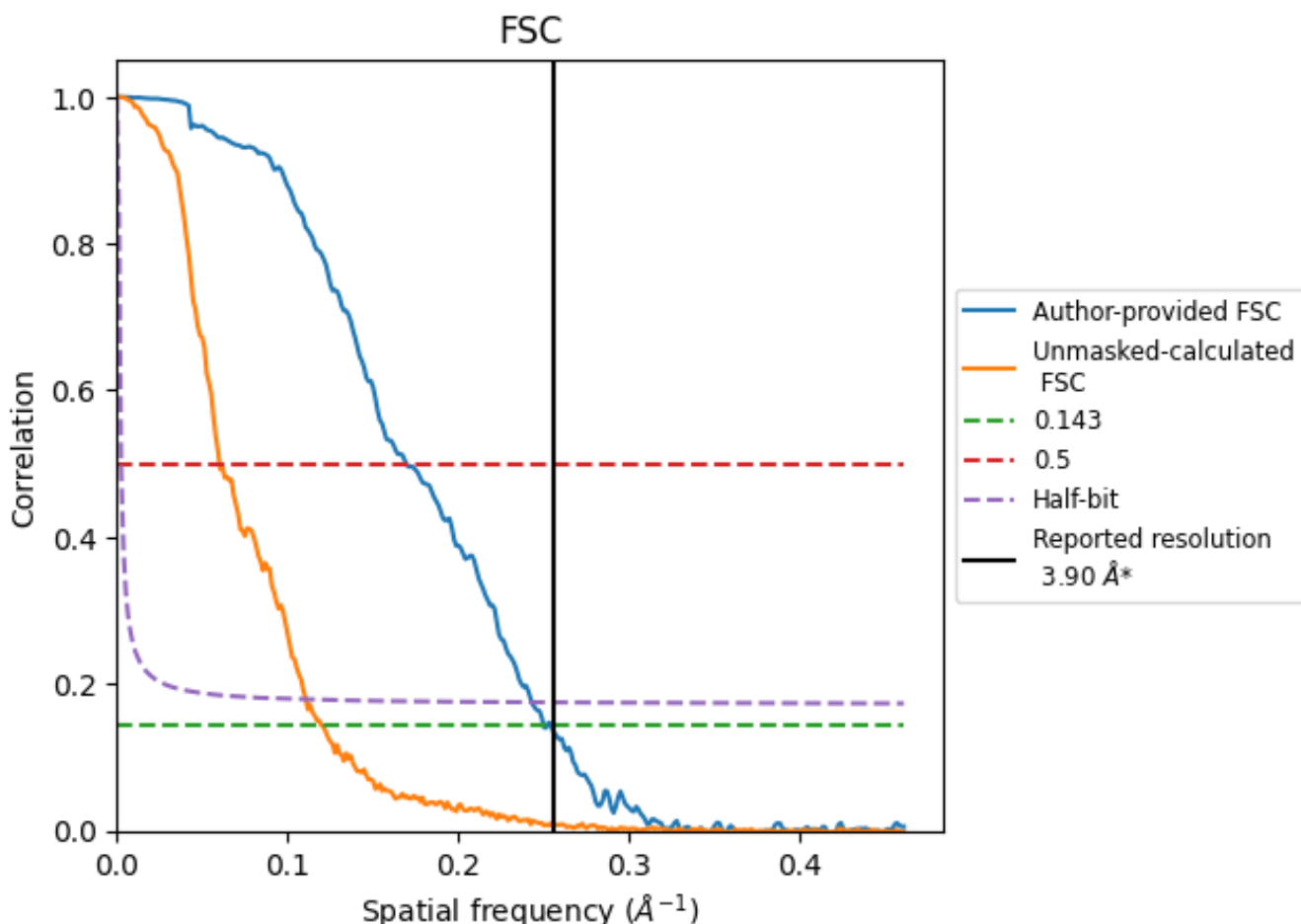


*Reported resolution corresponds to spatial frequency of 0.256 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.256 Å⁻¹

8.2 Resolution estimates [i](#)

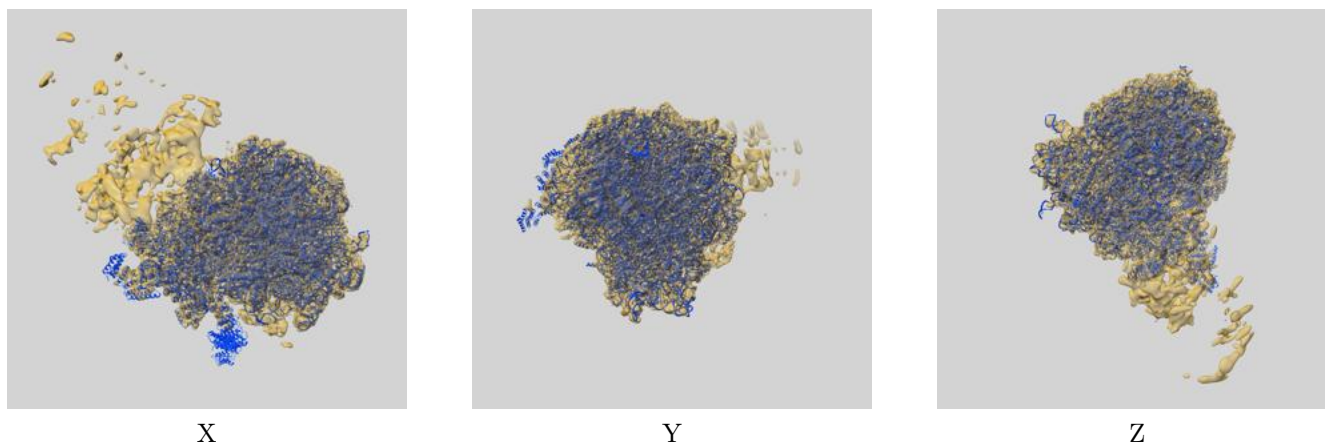
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.90	-	-
Author-provided FSC curve	4.00	5.90	4.11
Unmasked-calculated*	8.33	16.61	9.03

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 8.33 differs from the reported value 3.9 by more than 10 %

9 Map-model fit [i](#)

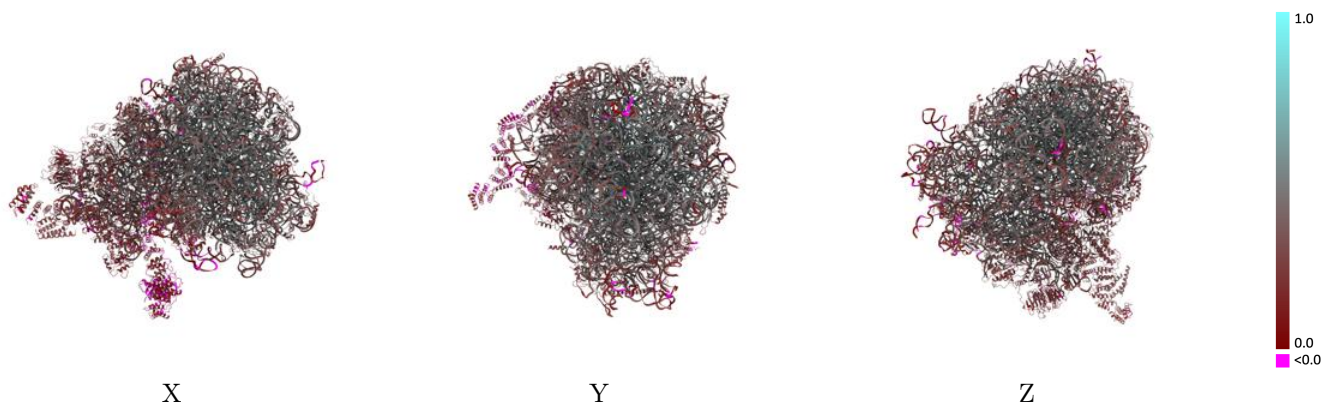
This section contains information regarding the fit between EMDB map EMD-12534 and PDB model 7NRC. Per-residue inclusion information can be found in section 3 on page 20.

9.1 Map-model overlay [i](#)



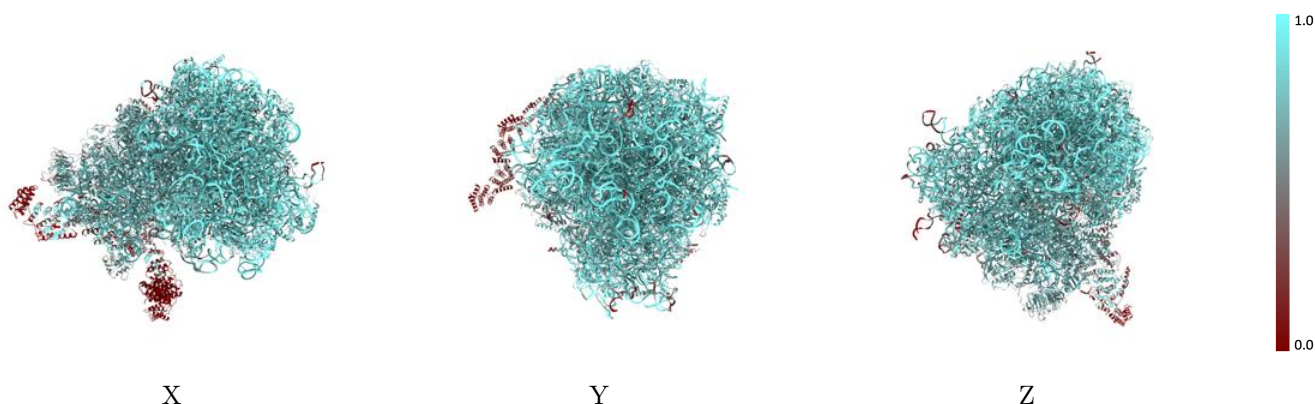
The images above show the 3D surface view of the map at the recommended contour level 0.03 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



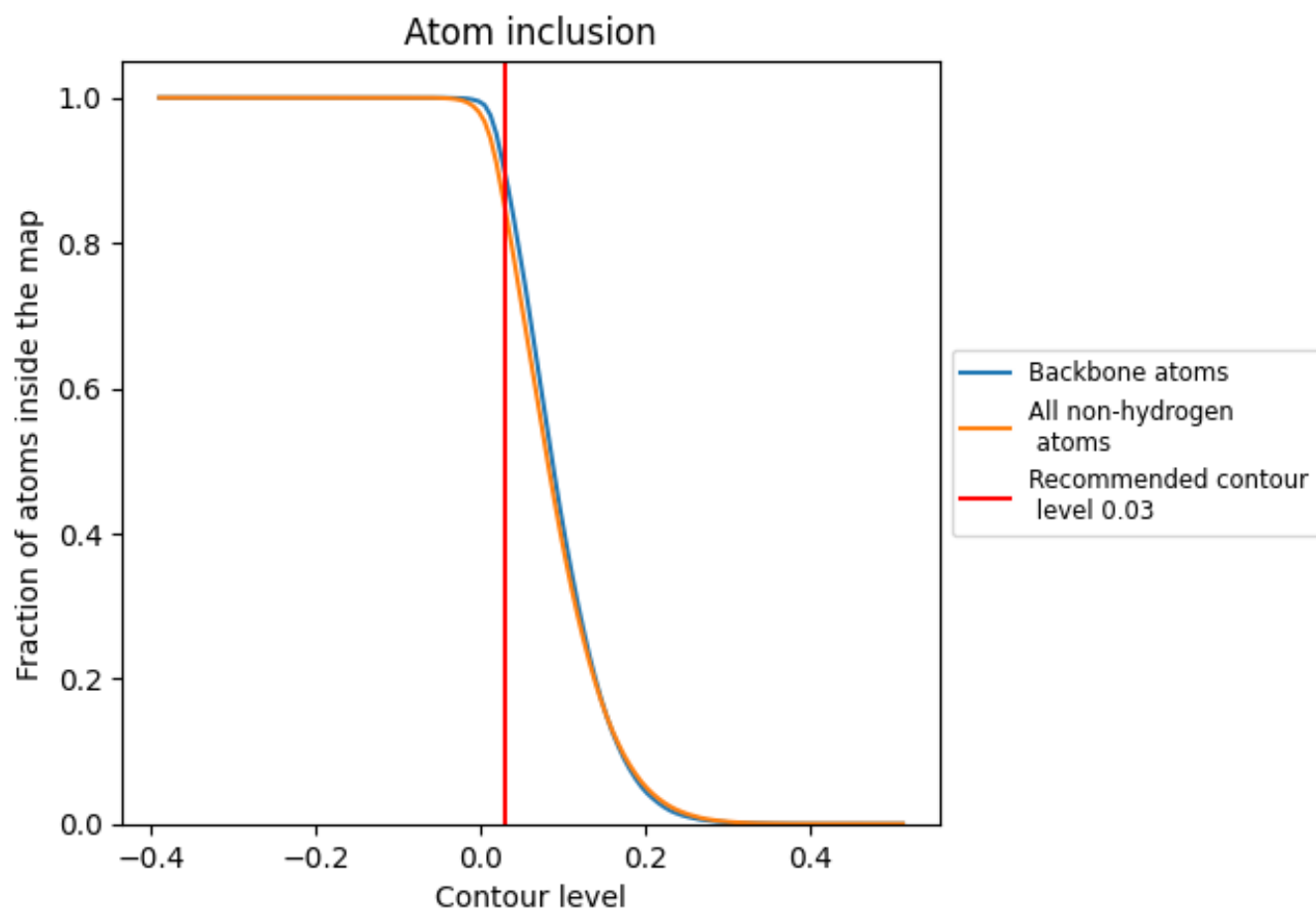
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.03).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 89% of all backbone atoms, 85% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary





















































































The table lists the average atom inclusion at the recommended contour level (0.03) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8451	 0.3780
A	 0.2817	 0.1550
LA	 0.9447	 0.4340
LB	 0.9492	 0.3610
LC	 0.9523	 0.4520
LD	 0.8641	 0.4660
LE	 0.8642	 0.4320
LF	 0.8481	 0.4070
LG	 0.7675	 0.2890
LH	 0.8259	 0.3730
LI	 0.8668	 0.4010
LJ	 0.8087	 0.3590
LK	 0.8108	 0.3720
LL	 0.7858	 0.3500
LM	 0.6147	 0.1650
LN	 0.8406	 0.3960
LO	 0.8606	 0.3760
LP	 0.8803	 0.4560
LQ	 0.8738	 0.4510
LR	 0.8628	 0.4320
LS	 0.8552	 0.4130
LT	 0.7835	 0.3720
LU	 0.8393	 0.4170
LV	 0.8430	 0.4050
LW	 0.7864	 0.3220
LX	 0.8212	 0.4320
LY	 0.8496	 0.4500
LZ	 0.8323	 0.4090
La	 0.8349	 0.3670
Lb	 0.8310	 0.3800
Lc	 0.8656	 0.4240
Ld	 0.8518	 0.3950
Le	 0.8041	 0.3710
Lf	 0.8118	 0.4100
Lg	 0.8668	 0.4440























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Chain	Atom inclusion	Q-score
Lh	 0.8843	 0.4680
Li	 0.8061	 0.4110
Lj	 0.8218	 0.3790
Lk	 0.8095	 0.3610
Ll	 0.8950	 0.4670
Lm	 0.7863	 0.3300
Ln	 0.8747	 0.4410
Lo	 0.8258	 0.3990
Lp	 0.8029	 0.4130
Lq	 0.8348	 0.4320
Lr	 0.8281	 0.4270
Ls	 0.6434	 0.3290
Lt	 0.8381	 0.2970
S2	 0.9019	 0.3700
SA	 0.6954	 0.3000
SB	 0.6816	 0.2860
SC	 0.6762	 0.2400
SD	 0.5780	 0.1730
SE	 0.7393	 0.2740
SF	 0.7312	 0.3150
SG	 0.6800	 0.2890
SH	 0.7009	 0.2750
SI	 0.6772	 0.2650
SJ	 0.7046	 0.2900
SK	 0.6076	 0.2620
SL	 0.6956	 0.3260
SM	 0.7849	 0.3670
SN	 0.5746	 0.1420
SO	 0.6183	 0.2390
SP	 0.7104	 0.3030
SQ	 0.6704	 0.3000
SR	 0.7780	 0.3720
SS	 0.7967	 0.3580
ST	 0.7682	 0.3090
SU	 0.6482	 0.2620
SV	 0.8285	 0.3660
SW	 0.7495	 0.3100
SX	 0.7894	 0.3990
SY	 0.7628	 0.3580
SZ	 0.7424	 0.3590
Sa	 0.7412	 0.3490
Sb	 0.8136	 0.3970

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Chain	Atom inclusion	Q-score
Sc	 0.8090	 0.4200
Sd	 0.7744	 0.3140
Se	 0.8136	 0.3980
Sf	 0.7388	 0.3380
Sg	 0.7494	 0.3190
Sl	 0.9466	 0.4400
Sm	 0.8870	 0.3140
Sn	 0.8954	 0.3330
So	 0.5370	 0.2640
Sp	 0.7583	 0.2700